

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050995.D
 Acq On : 12 Nov 2021 10:58
 Operator : CG/JU
 Sample : M4615-05
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0V04

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Title : SVOA CALIBRATION

Signal : TIC: BG050995.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.526	3	6	26	rVB	6996448	12074258	10.43%	2.760%
2	5.276	298	304	310	rBV2	120191	233397	0.20%	0.053%
3	5.629	359	364	374	rVB2	365064	811622	0.70%	0.186%
4	5.764	380	387	394	rVB	152405	269593	0.23%	0.062%
5	6.005	418	428	433	rBV	230884	413819	0.36%	0.095%
6	6.070	433	439	444	rVV2	462105	802580	0.69%	0.183%
7	6.158	450	454	458	rVV	406014	667742	0.58%	0.153%
8	6.211	458	463	472	rVB2	1363615	2421789	2.09%	0.554%
9	6.481	500	509	522	rBV5	504310	1894363	1.64%	0.433%
10	6.587	522	527	535	rVB2	207653	412927	0.36%	0.094%
11	6.710	535	548	551	rBV4	235129	793039	0.68%	0.181%
12	6.757	551	556	560	rVV	824278	1280212	1.11%	0.293%
13	6.810	560	565	566	rVV2	187827	252701	0.22%	0.058%
14	6.851	566	572	580	rVB4	1102779	2580167	2.23%	0.590%
15	6.963	586	591	597	rVB2	382235	625281	0.54%	0.143%
16	7.033	597	603	607	rBV3	196501	386406	0.33%	0.088%
17	7.098	607	614	619	rVV2	332470	819640	0.71%	0.187%
18	7.157	619	624	626	rVV2	293954	464183	0.40%	0.106%
19	7.198	626	631	635	rVV	1101859	1906647	1.65%	0.436%
20	7.251	635	640	645	rVV	1058203	1645193	1.42%	0.376%
21	7.309	645	650	654	rVV3	471693	826941	0.71%	0.189%
22	7.362	654	659	664	rVV2	1667229	3303377	2.85%	0.755%
23	7.632	664	705	709	rVV2	12051648	115795749	100.00%	26.473%
24	7.674	709	712	715	rVV2	629197	1168545	1.01%	0.267%
25	7.715	715	719	722	rVV2	1029770	1993817	1.72%	0.456%
26	7.750	722	725	729	rVV	1760399	2760495	2.38%	0.631%
27	7.791	729	732	735	rVV	592206	1030961	0.89%	0.236%
28	7.856	735	743	747	rVV	5482392	10376109	8.96%	2.372%
29	7.915	747	753	757	rVV2	771099	1797300	1.55%	0.411%
30	7.956	757	760	765	rVV2	532493	888789	0.77%	0.203%
31	8.008	765	769	773	rVV3	294032	582075	0.50%	0.133%
32	8.050	773	776	780	rVV3	299139	497260	0.43%	0.114%
33	8.144	780	792	795	rVV5	449396	1455348	1.26%	0.333%
34	8.197	796	801	809	rVV	1933544	3485299	3.01%	0.797%
35	8.285	809	816	823	rVV4	435519	1081190	0.93%	0.247%
36	8.367	823	830	834	rVV3	543733	1336125	1.15%	0.305%

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37	8.408	834	837	844	rVV4	574409	1404449	1.21%	0.321%
38	8.473	844	848	851	rVV2	549053	904580	0.78%	0.207%
39	8.520	851	856	861	rVV2	1483943	2693928	2.33%	0.616%
40	8.573	861	865	871	rVB2	350130	653201	0.56%	0.149%

41	8.643	871	877	879	rBV5	157944	288837	0.25%	0.066%
42	8.672	879	882	887	rVV3	212672	344875	0.30%	0.079%
43	8.743	887	894	897	rVV2	717268	1430848	1.24%	0.327%
44	8.790	897	902	911	rVB4	1071350	3002749	2.59%	0.686%
45	8.878	911	917	924	rBV3	1208028	2120693	1.83%	0.485%

46	8.984	929	935	943	rVV2	762250	1885880	1.63%	0.431%
47	9.084	947	952	966	rVB3	669855	1775099	1.53%	0.406%
48	9.248	977	980	985	rVB	300074	465965	0.40%	0.107%
49	9.301	985	989	991	rBV2	285142	422680	0.37%	0.097%
50	9.342	991	996	1006	rVV3	657252	1608923	1.39%	0.368%

51	9.454	1006	1015	1020	rVV	3665171	7057869	6.10%	1.614%
52	9.554	1026	1032	1034	rVV4	222027	433099	0.37%	0.099%
53	9.595	1035	1039	1046	rVB4	337113	686038	0.59%	0.157%
54	9.695	1046	1056	1064	rBV5	259683	661174	0.57%	0.151%
55	9.795	1066	1073	1079	rBV6	89985	279400	0.24%	0.064%

56	9.865	1080	1085	1092	rVB3	302959	594514	0.51%	0.136%
57	9.930	1092	1096	1100	rBV	185760	318254	0.27%	0.073%
58	9.971	1100	1103	1117	rVB4	189933	420912	0.36%	0.096%
59	10.159	1127	1135	1143	rVB3	331967	661648	0.57%	0.151%
60	10.229	1143	1147	1151	rBV3	124847	219799	0.19%	0.050%

61	10.447	1178	1184	1197	rVV6	241282	581320	0.50%	0.133%
62	10.993	1269	1277	1284	rBV2	472244	823322	0.71%	0.188%
63	11.064	1284	1289	1294	rVV	180737	330429	0.29%	0.076%
64	11.634	1375	1386	1401	rBV	1846938	3826048	3.30%	0.875%
65	11.851	1407	1423	1438	rBV2	3640624	16549550	14.29%	3.783%

66	12.074	1444	1461	1485	rVB3	761209	5358146	4.63%	1.225%
67	12.397	1506	1516	1526	rBV	710564	1266787	1.09%	0.290%
68	13.244	1645	1660	1677	rVB	1114878	2759841	2.38%	0.631%
69	13.561	1701	1714	1732	rBV	1038418	2737928	2.36%	0.626%
70	14.554	1877	1883	1894	rVB	160868	271211	0.23%	0.062%

71	14.859	1927	1935	1944	rVB	302442	498513	0.43%	0.114%
72	15.841	2094	2102	2115	rBV2	1055954	1656803	1.43%	0.379%
73	17.603	2396	2402	2408	rVB	335648	530573	0.46%	0.121%
74	17.697	2412	2418	2427	rBV	202175	325439	0.28%	0.074%
75	18.214	2502	2506	2512	rVB	173565	273204	0.24%	0.062%

76	18.473	2544	2550	2559	rBV	300943	486790	0.42%	0.111%
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77	18.772	2592	2601	2610	rVV	3693185	6461017	5.58%	1.477%
78	19.060	2630	2650	2672	rVB	13969430	54980880	47.48%	12.569%
79	19.319	2688	2694	2700	rBV2	735387	1373649	1.19%	0.314%
80	19.477	2702	2721	2729	rVV	13601226	50402867	43.53%	11.523%
81	19.642	2741	2749	2760	rVV	344463	799005	0.69%	0.183%
82	19.795	2770	2775	2786	rVB2	212792	534120	0.46%	0.122%
83	19.983	2802	2807	2811	rBV	263070	411957	0.36%	0.094%
84	20.147	2832	2835	2844	rVB3	148287	233477	0.20%	0.053%
85	20.494	2889	2894	2902	rVB2	1187645	1657658	1.43%	0.379%
86	20.570	2902	2907	2917	rBV2	264559	849519	0.73%	0.194%
87	20.653	2917	2921	2933	rVB2	215050	552203	0.48%	0.126%
88	20.823	2944	2950	2954	rBV3	180106	342673	0.30%	0.078%
89	20.923	2961	2967	2974	rBV5	105656	309150	0.27%	0.071%
90	21.557	3050	3075	3080	rBV3	13346305	58671677	50.67%	13.413%
91	21.692	3094	3098	3104	rVB	518664	1079055	0.93%	0.247%
92	21.910	3130	3135	3143	rVB	267189	559994	0.48%	0.128%
93	22.116	3164	3170	3184	rBV	476350	1386875	1.20%	0.317%
94	22.738	3271	3276	3280	rBV2	112408	260312	0.22%	0.060%
95	22.803	3281	3287	3295	rVV	783664	1740018	1.50%	0.398%
96	25.312	3707	3714	3724	rVB2	169028	524682	0.45%	0.120%
97	25.447	3730	3737	3747	rVB	92052	248910	0.21%	0.057%
98	26.463	3897	3910	3934	rBV2	559292	2482987	2.14%	0.568%
99	26.875	3970	3980	4002	rVB2	298663	1207594	1.04%	0.276%
100	27.544	4079	4094	4114	rBV3	551527	2600520	2.25%	0.595%

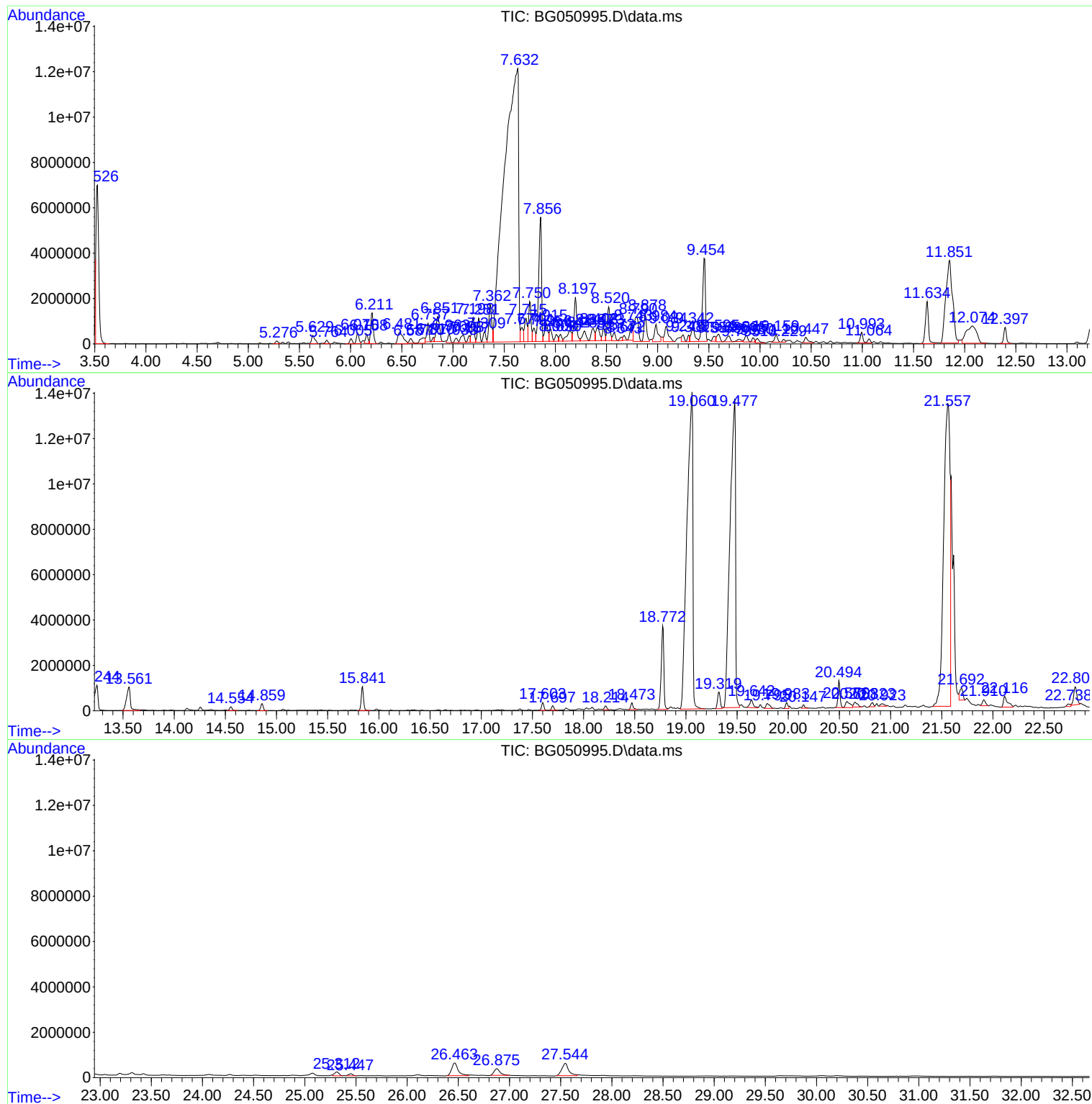
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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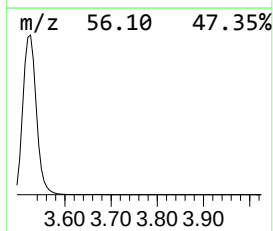
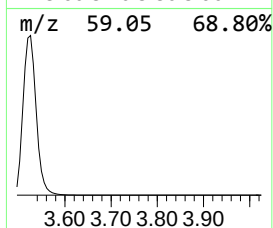
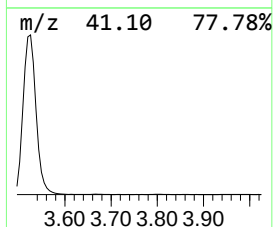
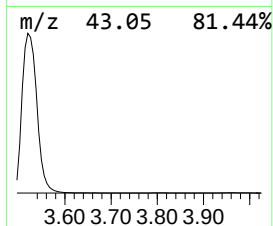
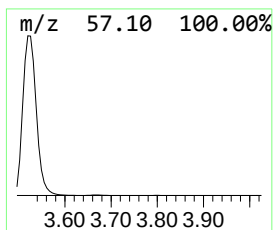
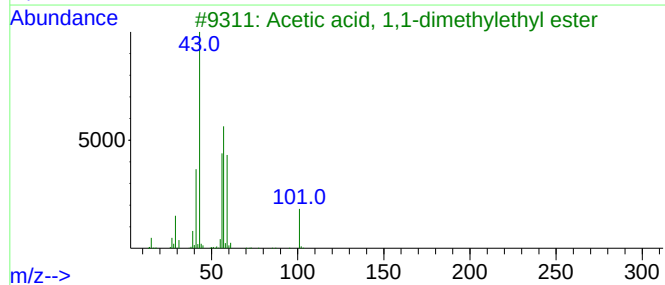
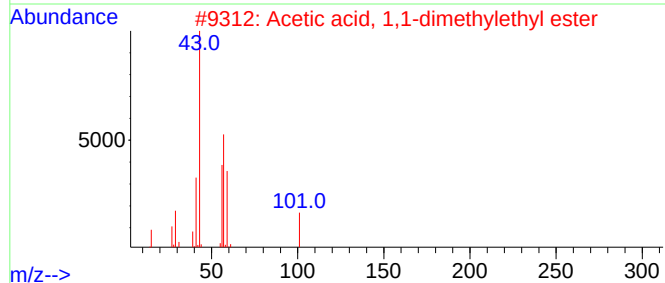
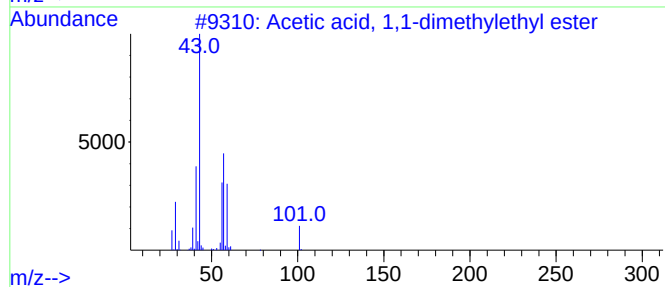
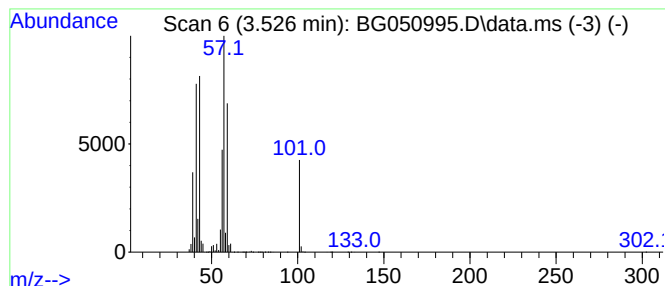
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Acetic acid, 1,1-dimethylet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.526	223.35 ng/ul	12074300	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	72		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	64		
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	64		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	64		
5	Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	38		



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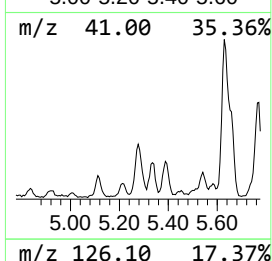
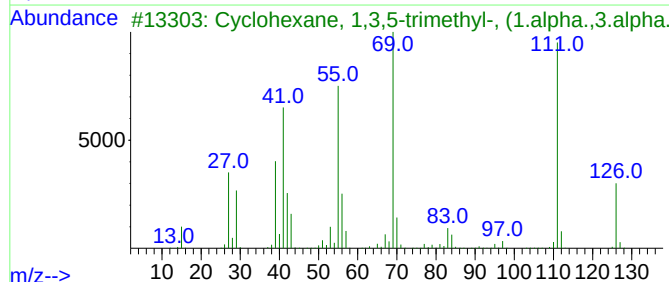
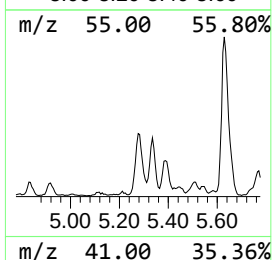
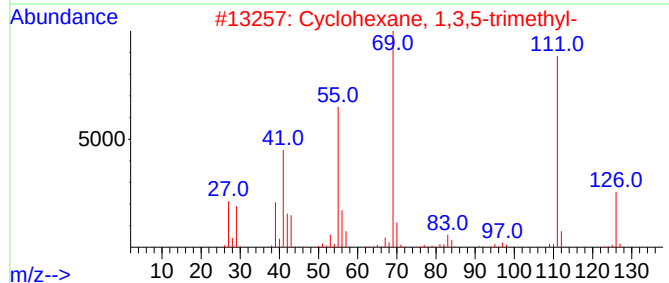
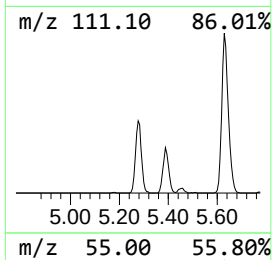
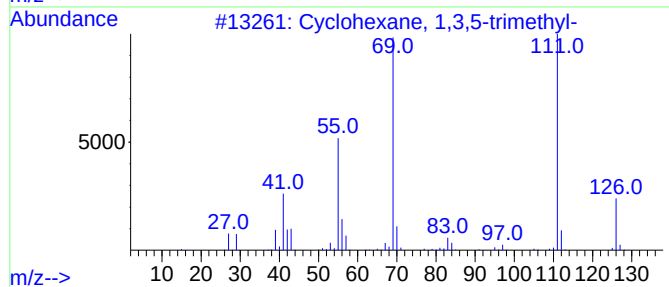
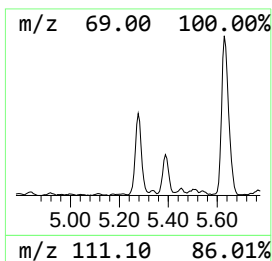
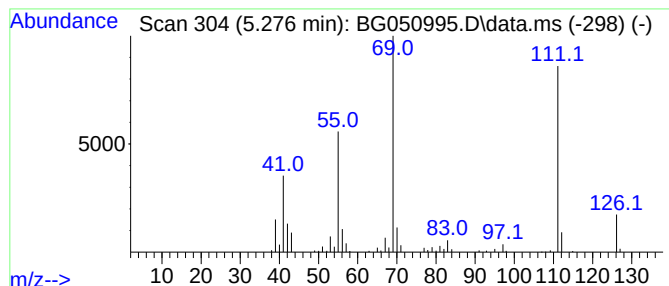
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Cyclic5.276 Concentration Rank 79

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.276	4.32 ng/ul	233397	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	91
2			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	90
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	87
4			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	76
5			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	76



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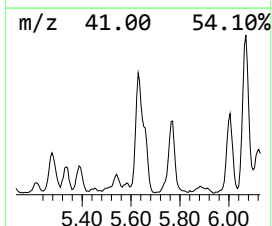
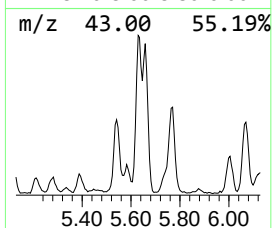
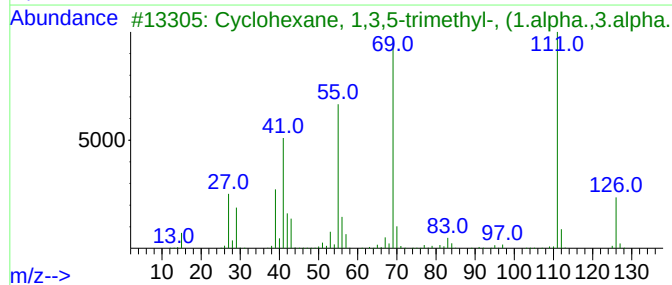
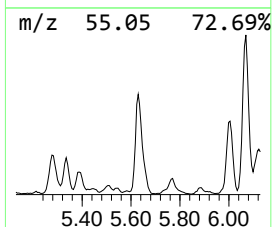
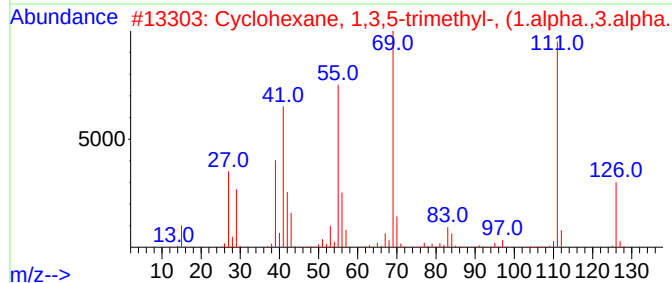
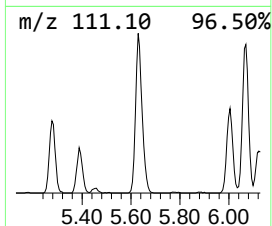
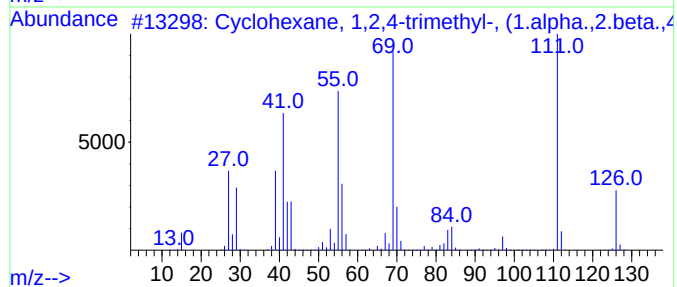
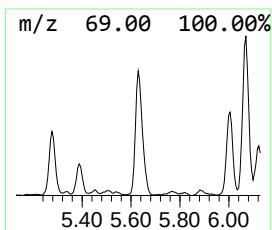
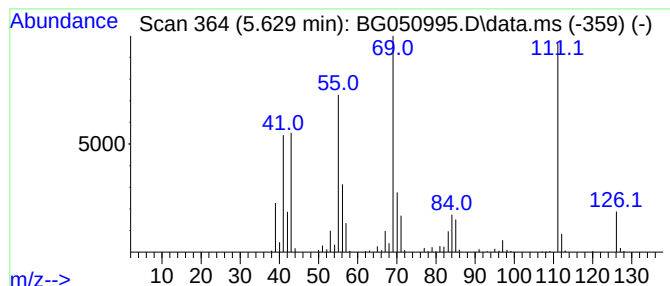
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Cyclic5.629 Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.629	15.01 ng/ul	811622	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	90
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	83
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	72
4			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	52
5			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
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Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

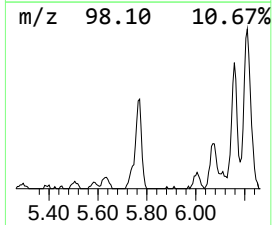
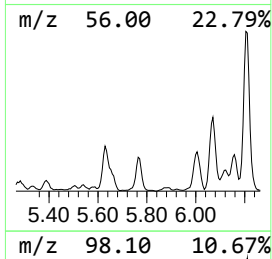
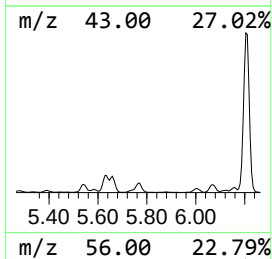
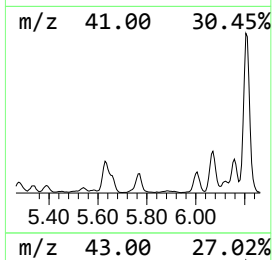
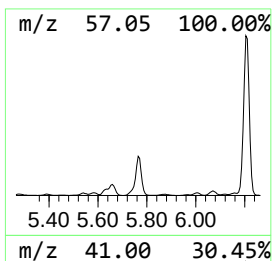
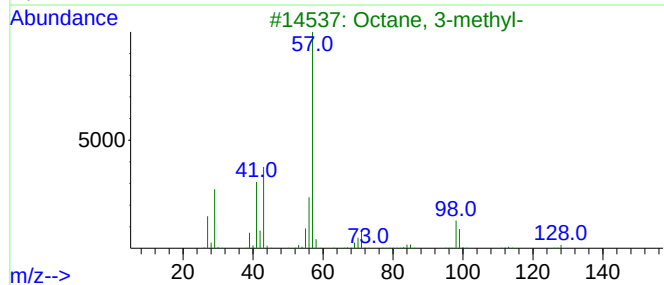
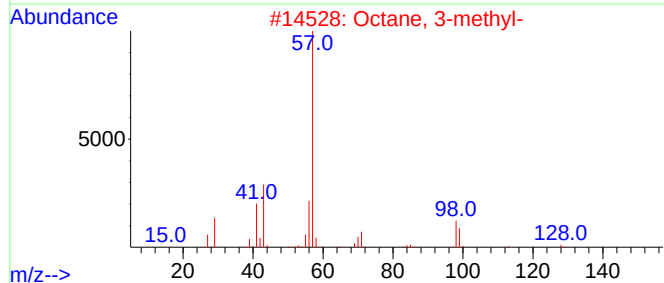
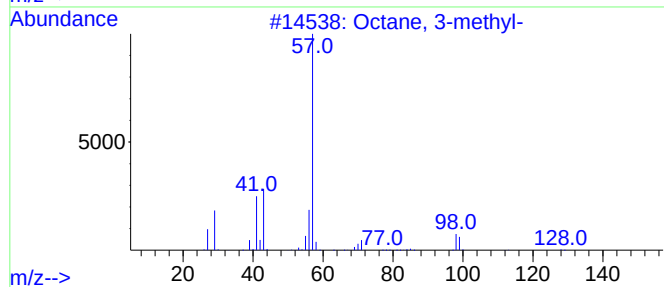
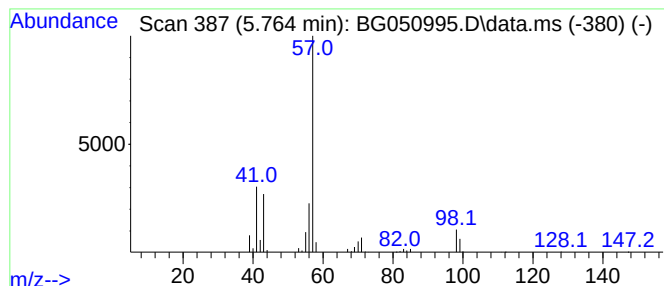
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 77

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.764	4.99 ng/ul	269593	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3-methyl-	128	C9H20	002216-33-3	90
2			Octane, 3-methyl-	128	C9H20	002216-33-3	87
3			Octane, 3-methyl-	128	C9H20	002216-33-3	74
4			Undecane, 6-methyl-	170	C12H26	017302-33-9	59
5			Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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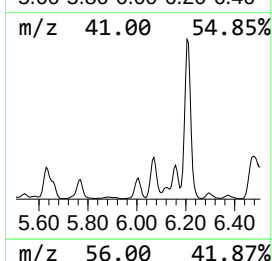
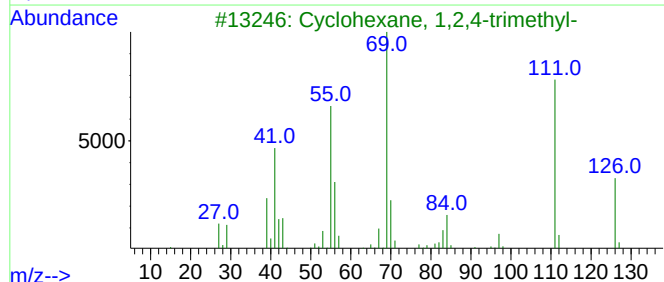
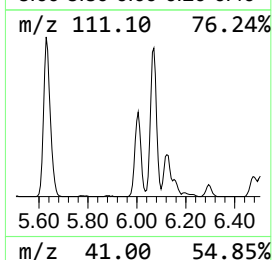
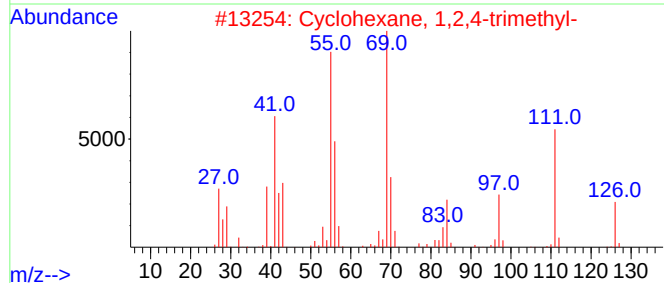
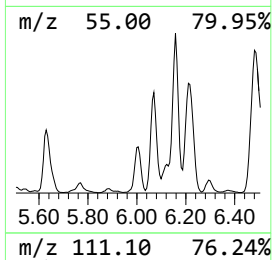
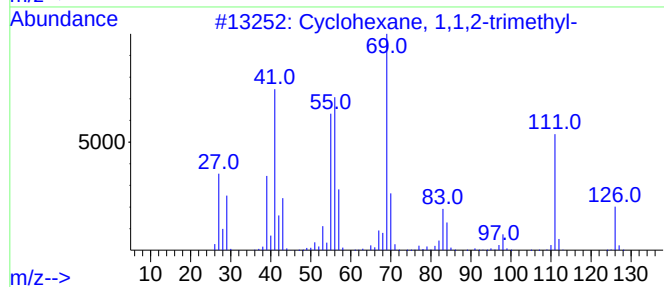
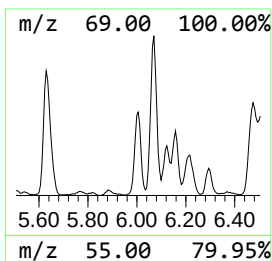
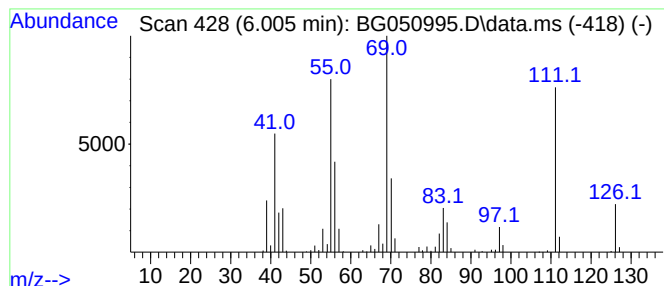
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic6.005 Concentration Rank 73

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.005	7.65 ng/ul	413819	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	90
2			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
3			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
4			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	81
5			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

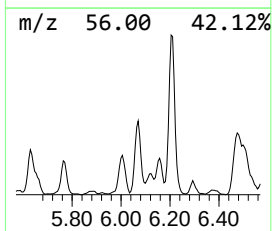
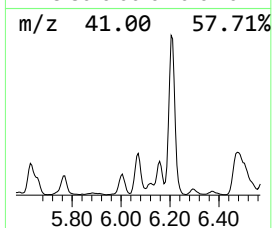
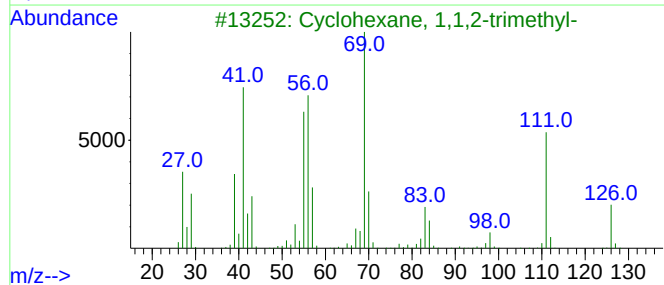
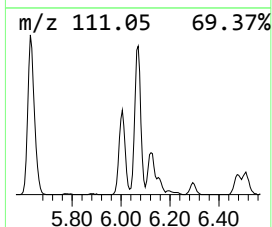
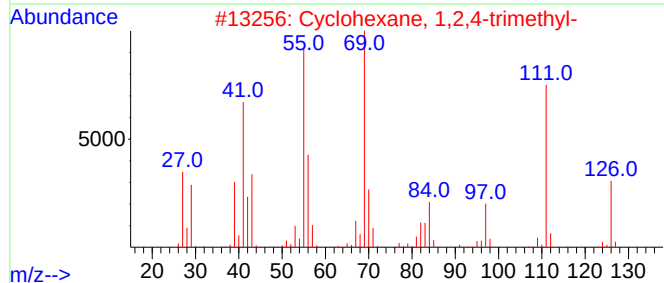
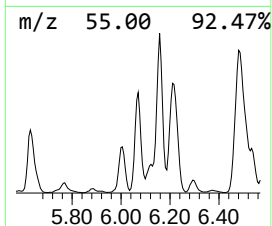
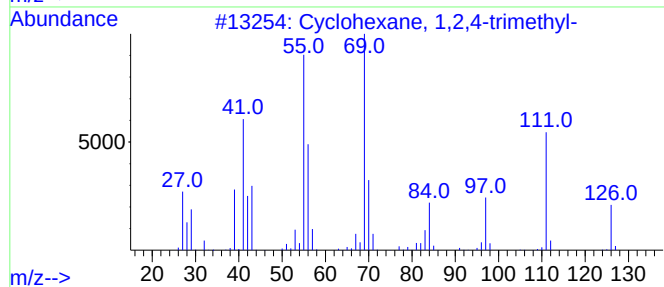
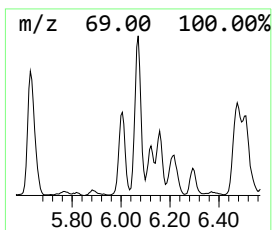
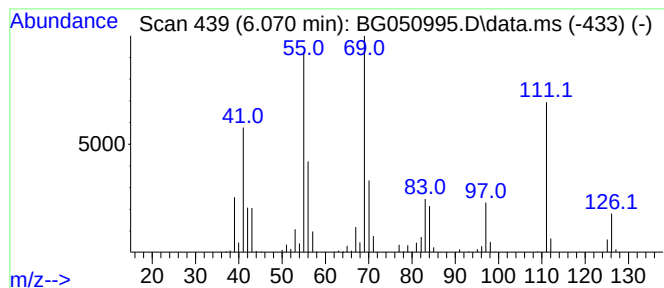
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic6.070 Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.070	14.85 ng/ul	802580	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	97
2			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
3			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	70
4			Cyclopentane, propyl-	112	C8H16	002040-96-2	70
5			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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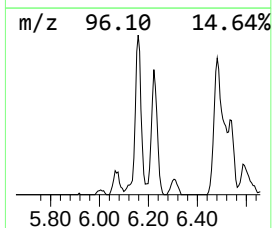
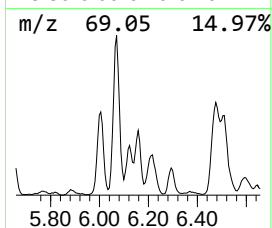
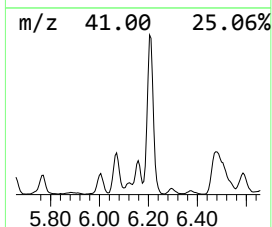
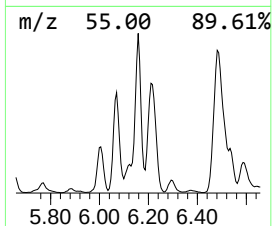
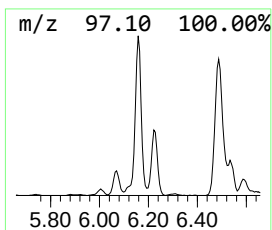
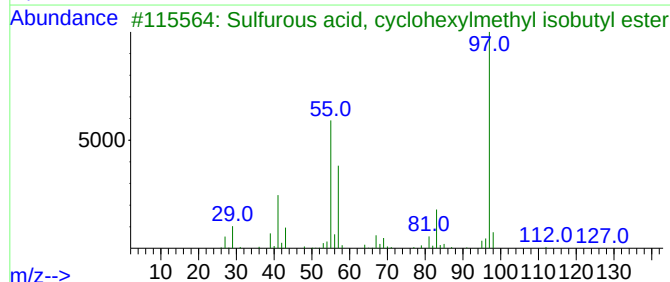
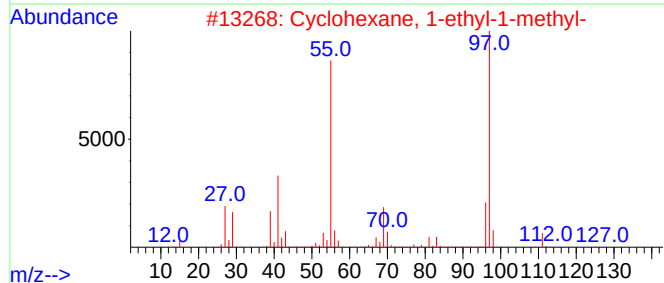
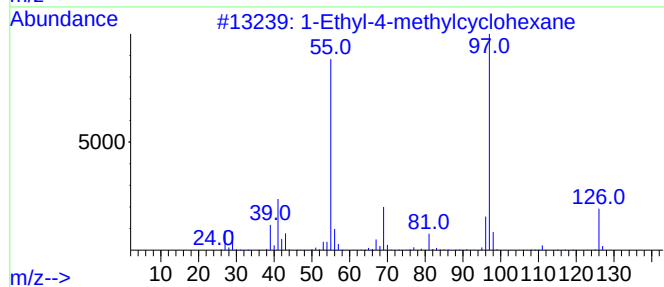
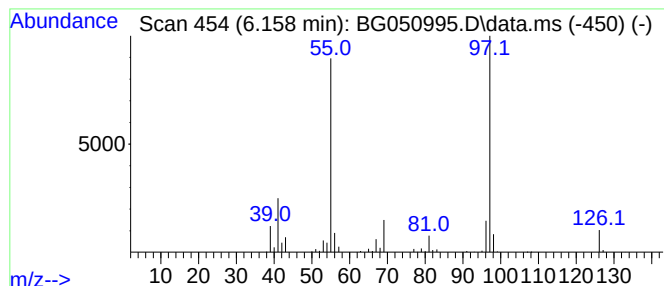
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.158	12.35 ng/ul	667742	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	87
2		Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	78
3		Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	72
4		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	72
5		Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
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Instrument :
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ClientSampleId :
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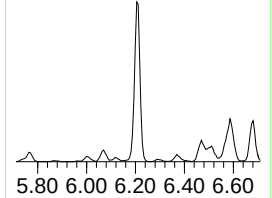
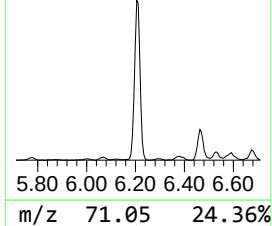
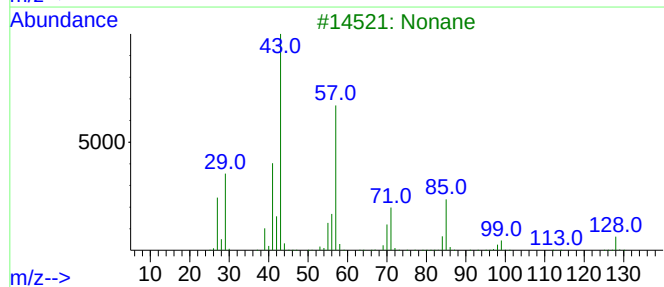
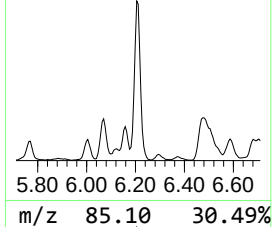
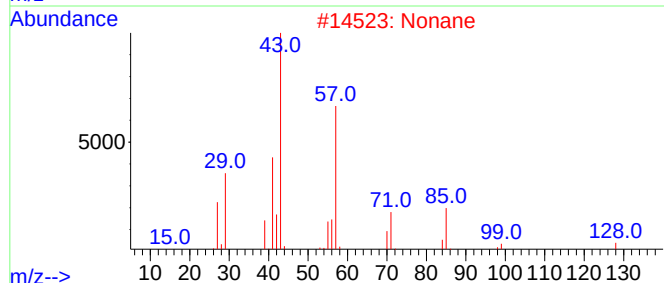
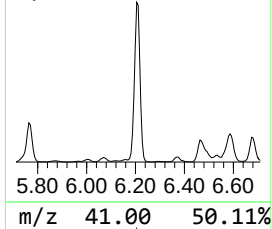
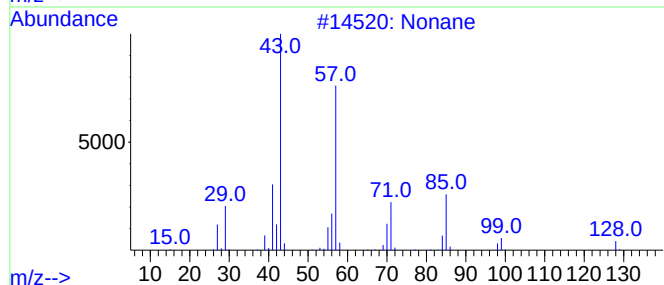
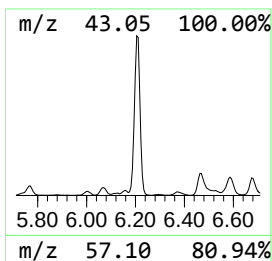
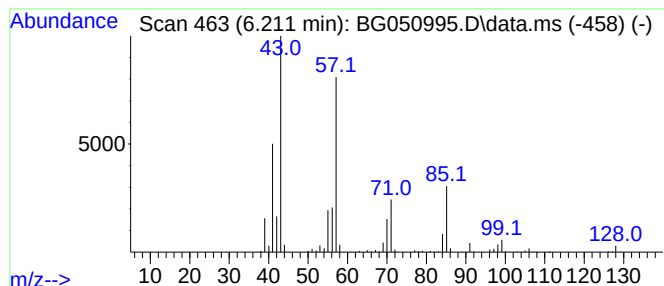
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.211	44.80 ng/ul	2421790	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	91
2			Nonane	128	C9H20	000111-84-2	91
3			Nonane	128	C9H20	000111-84-2	90
4			Octane	114	C8H18	000111-65-9	64
5			Octane	114	C8H18	000111-65-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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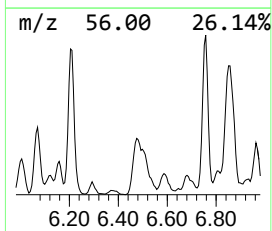
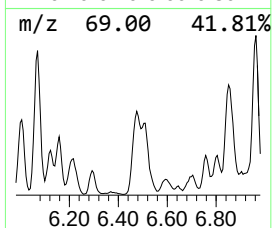
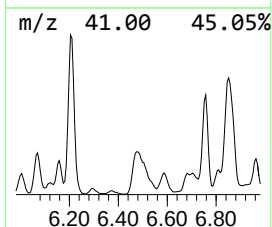
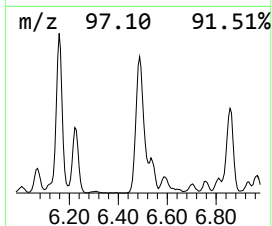
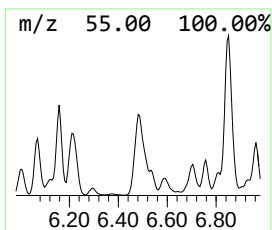
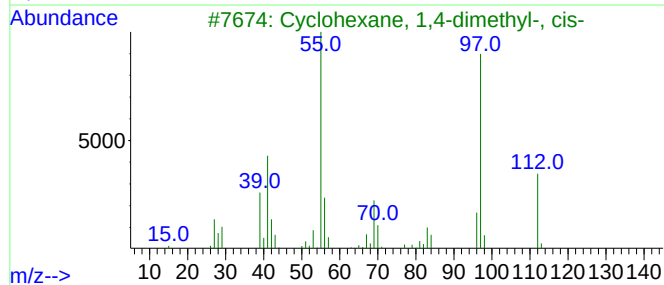
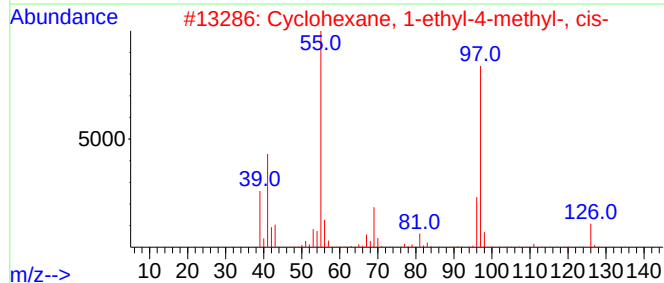
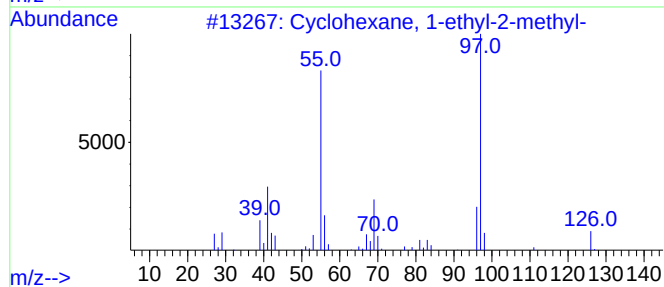
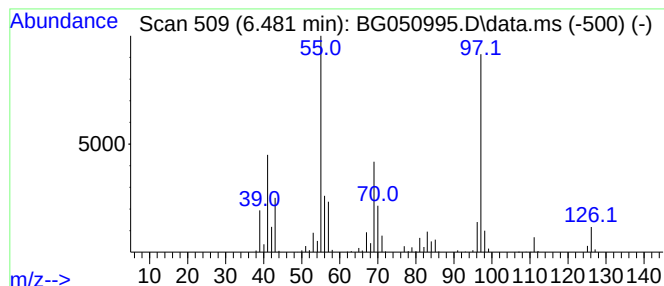
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Cyclic6.481 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.481	35.04 ng/ul	1894360	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	68
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	68
3			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	64
4			Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	64
5			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
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Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

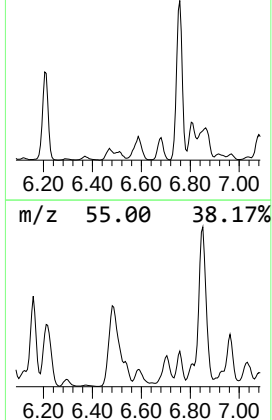
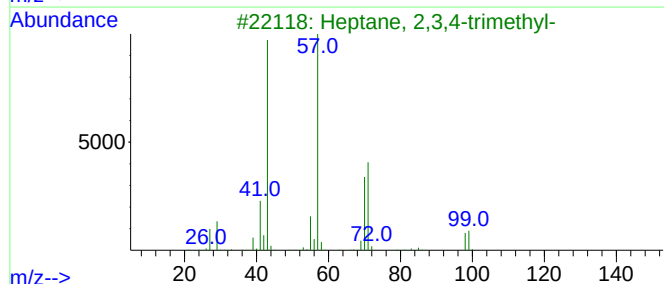
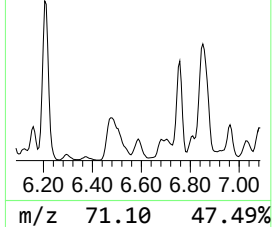
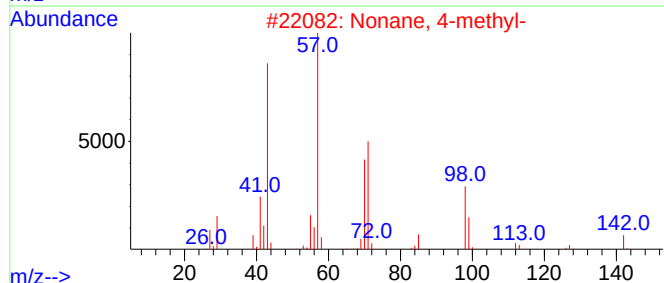
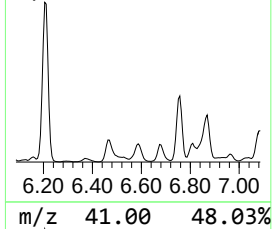
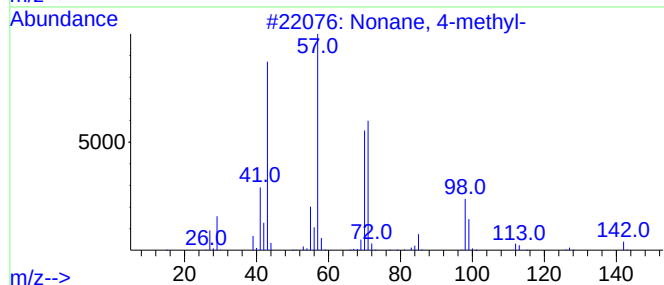
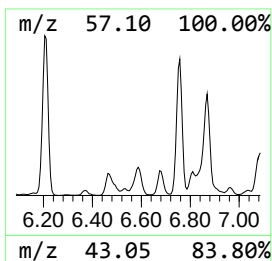
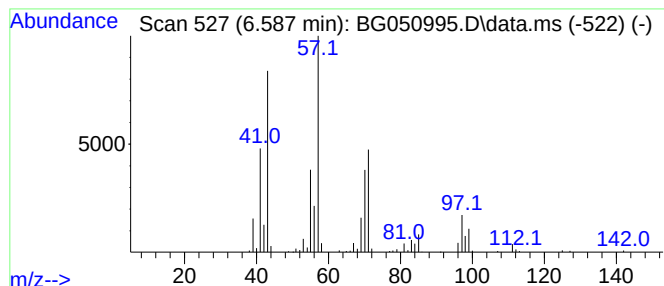
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 74

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.587	7.64 ng/ul	412927	1,4-Dichlorobenzene-d4			8.261
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-		142	C10H22	017301-94-9	59
2	Nonane, 4-methyl-		142	C10H22	017301-94-9	59
3	Heptane, 2,3,4-trimethyl-		142	C10H22	052896-95-4	59
4	Undecane, 4,5-dimethyl-		184	C13H28	017312-79-7	53
5	Nonane, 4-methyl-		142	C10H22	017301-94-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
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Instrument :
BNA_G
ClientSampleId :
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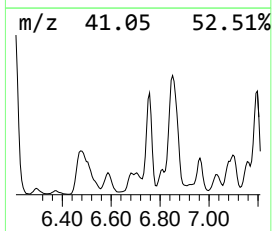
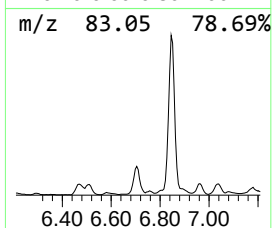
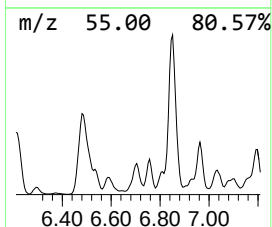
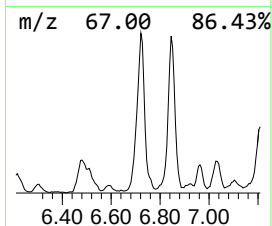
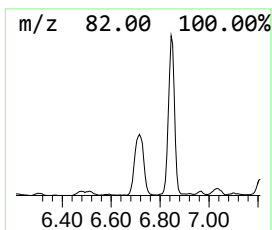
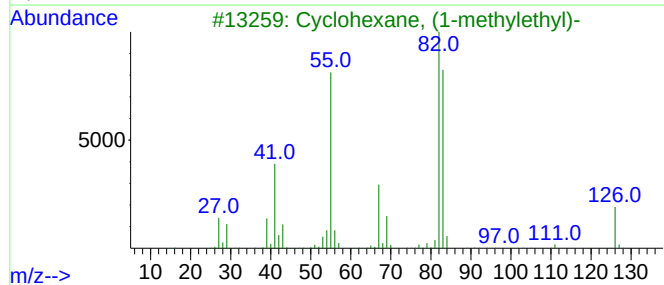
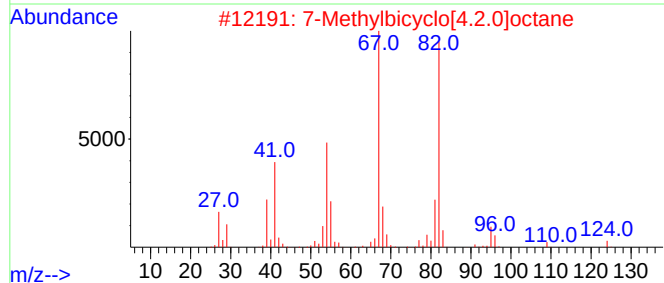
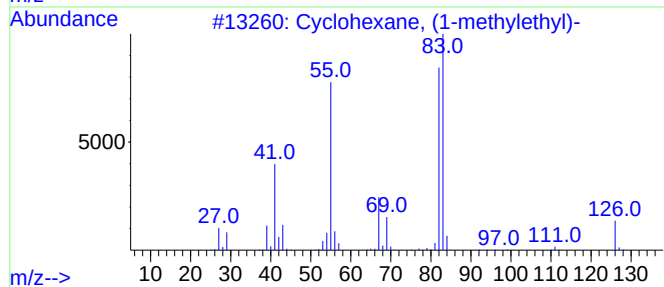
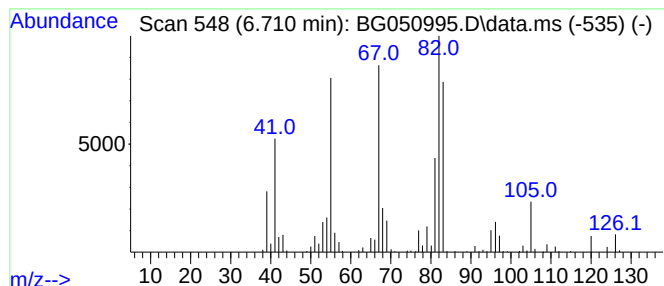
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Cyclic6.710 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.710	14.67 ng/ul	793039	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53
2			7-Methylbicyclo[4.2.0]octane	124	C9H16	1000210-90-2	53
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	47
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	43
5			Cyclopropane, (1-methylethylidene)-	82	C6H10	004741-86-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
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Instrument :
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ClientSampleId :
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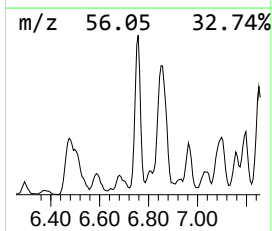
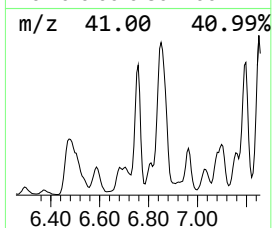
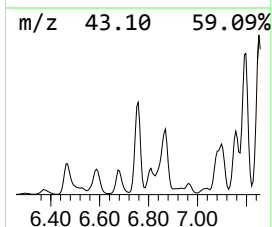
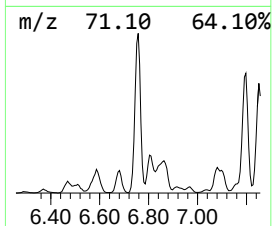
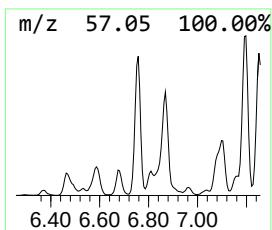
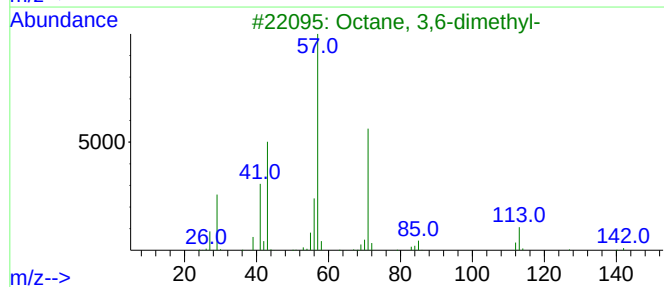
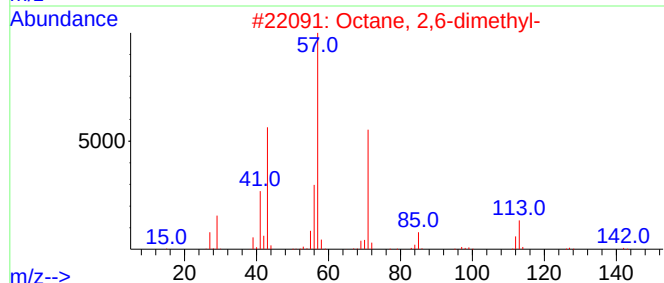
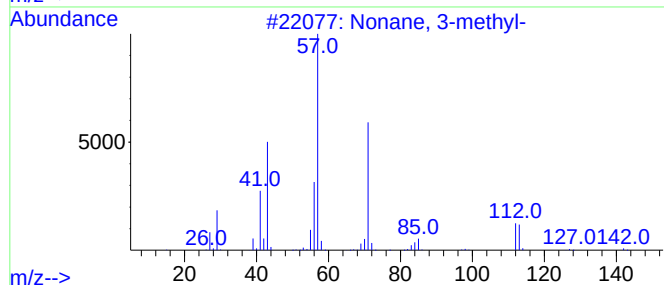
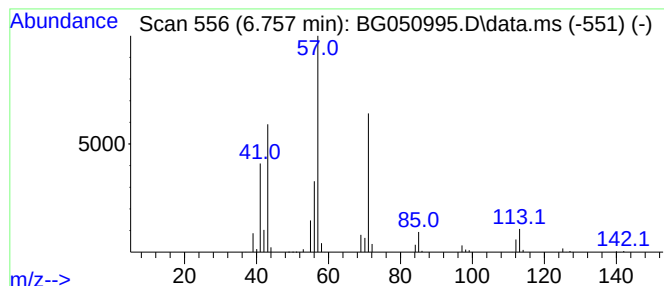
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.757	23.68 ng/ul	1280210	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	87
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	86
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4			Nonane, 3-methyl-	142	C10H22	005911-04-6	83
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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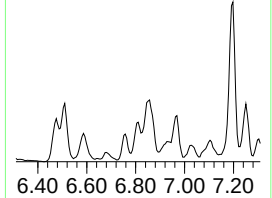
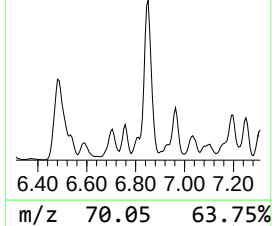
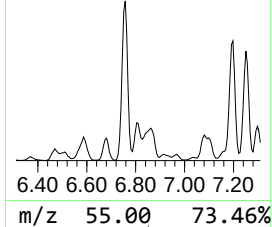
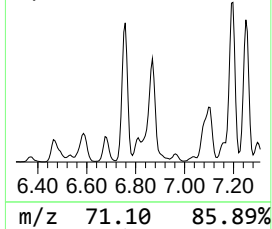
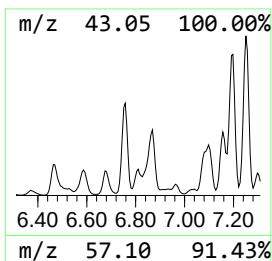
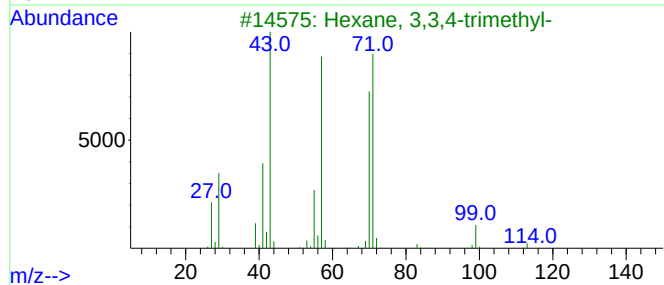
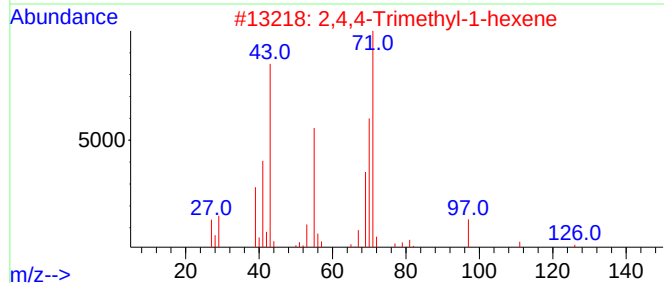
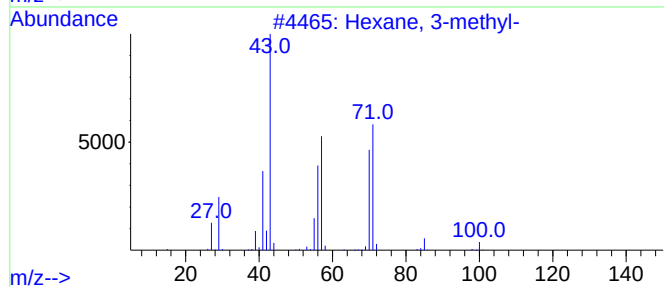
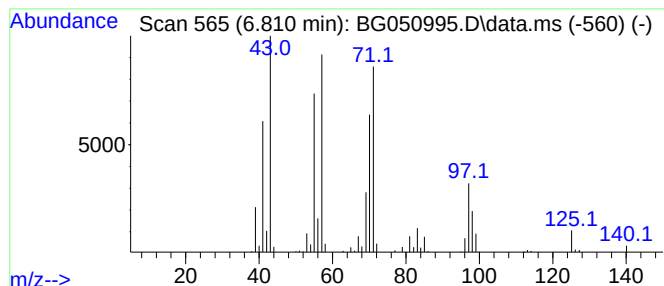
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 78

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.810	4.67 ng/ul	252701	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-			100	C7H16	000589-34-4	49
2	2,4,4-Trimethyl-1-hexene			126	C9H18	051174-12-0	49
3	Hexane, 3,3,4-trimethyl-			128	C9H20	016747-31-2	47
4	Hexane, 2,4,4-trimethyl-			128	C9H20	016747-30-1	47
5	Heptane, 3,3-dimethyl-			128	C9H20	004032-86-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
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ClientSampleId :
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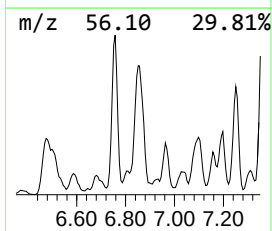
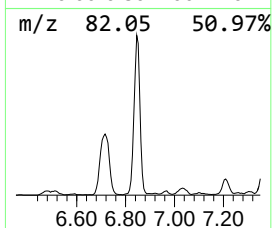
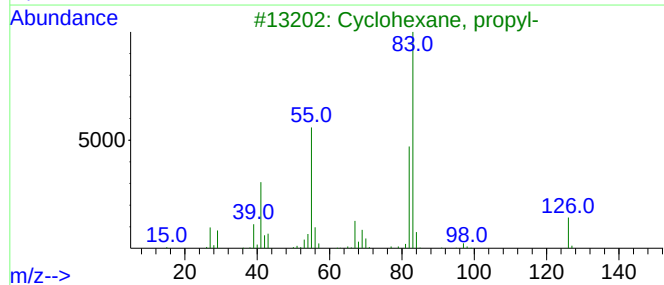
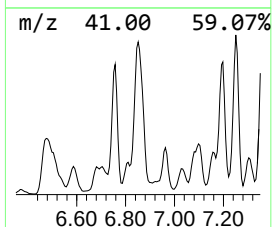
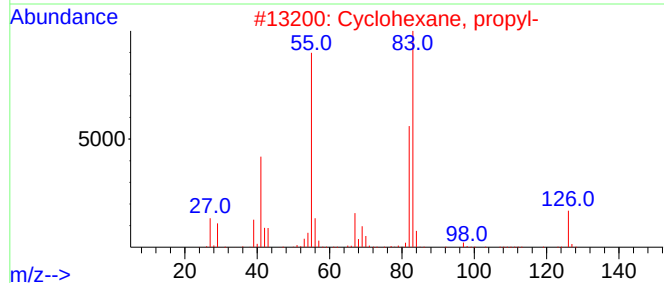
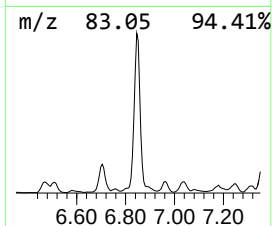
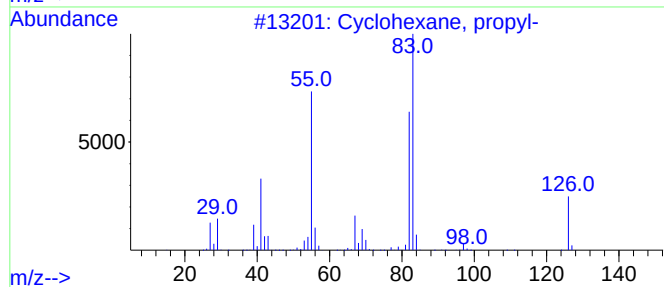
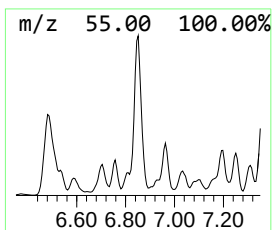
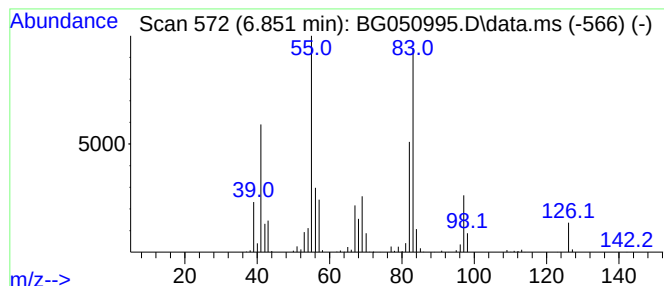
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Cyclic6.851 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.851	47.73 ng/ul	2580170	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	81
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
5			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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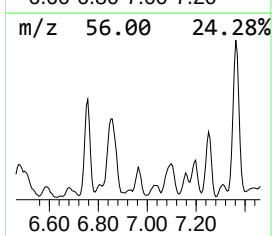
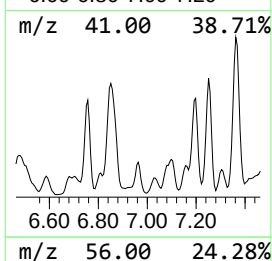
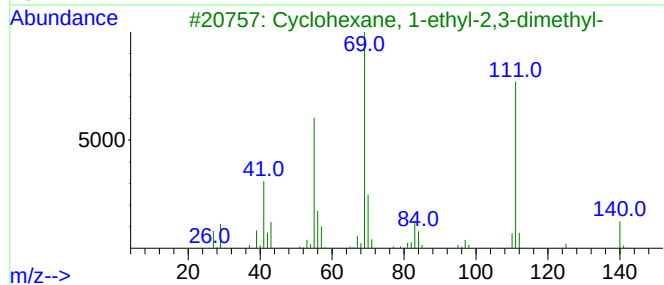
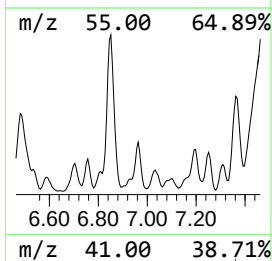
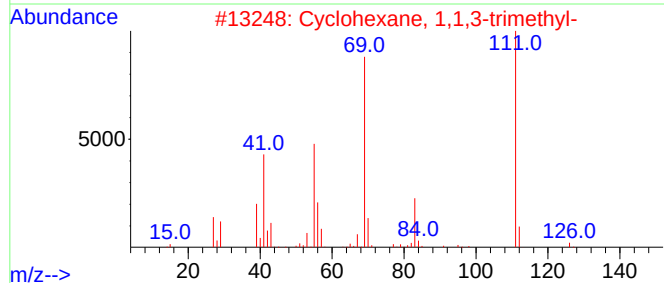
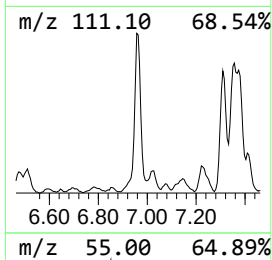
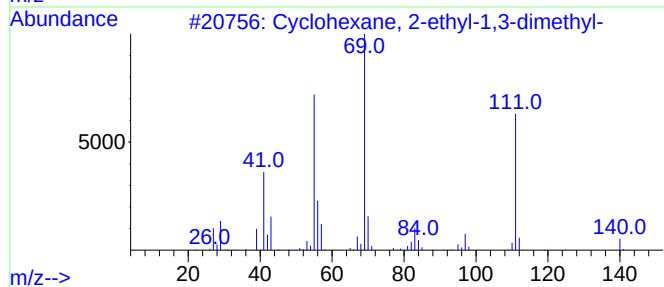
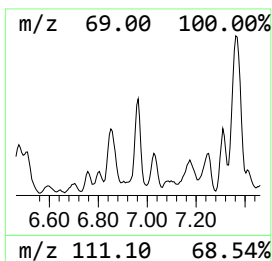
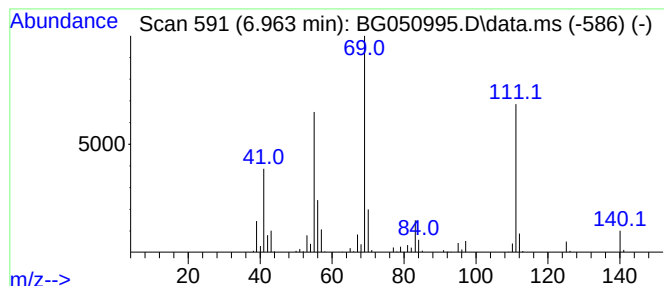
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 (DEL) Alkane: Cyclic6.963 Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.963	11.57 ng/ul	625281	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	93
2		Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	87
3		Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	87
4		Cyclooctane, butyl-	168	C12H24	016538-93-5	64
5		3-Hexene, 2,5-dimethyl-	112	C8H16	015910-22-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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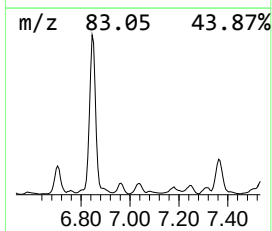
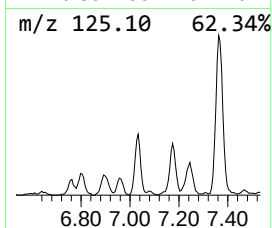
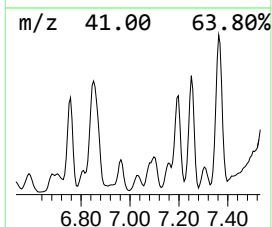
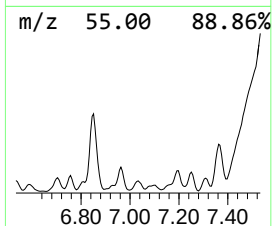
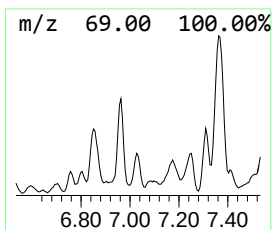
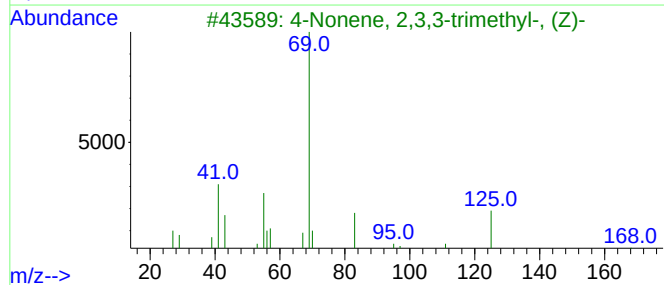
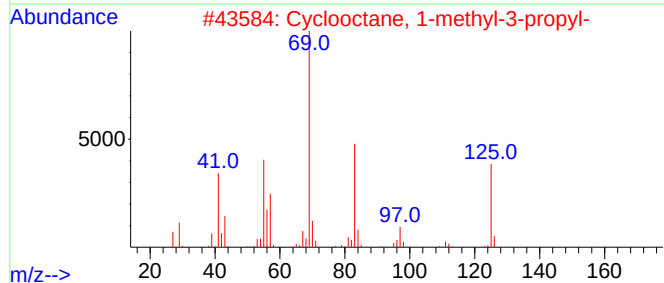
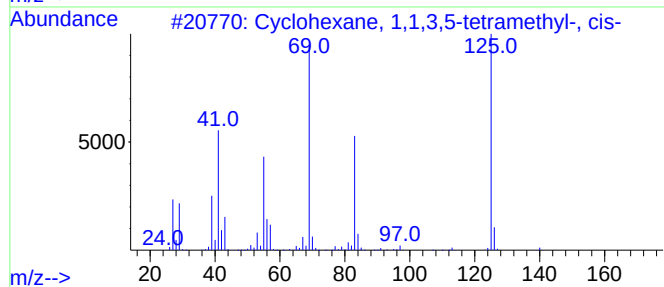
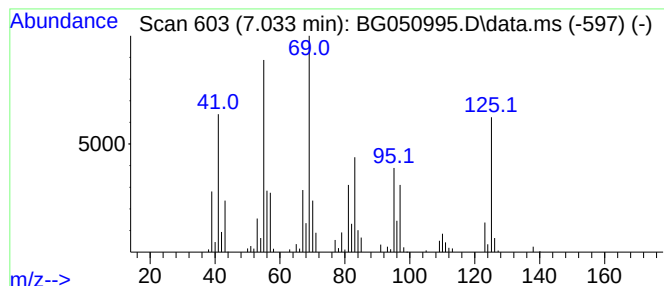
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Cyclic7.033 Concentration Rank 75

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.033	7.15 ng/ul	386406	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	47
2		Cyclooctane, 1-methyl-3-propyl-	168	C12H24	255885-37-1	43
3		4-Nonene, 2,3,3-trimethyl-, (Z)-	168	C12H24	063830-68-2	35
4		Cyclopentanecarboxylic acid, eth...	140	C8H12O2	016523-06-1	30
5		trans-7-Methyl-3-octene	126	C9H18	1000113-52-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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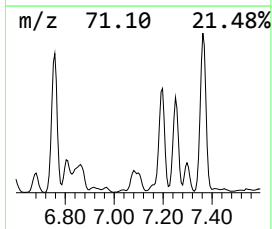
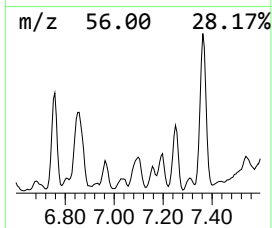
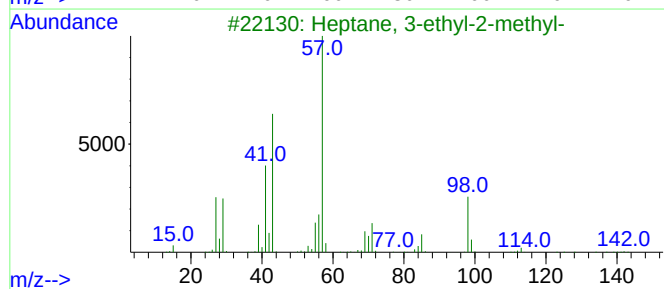
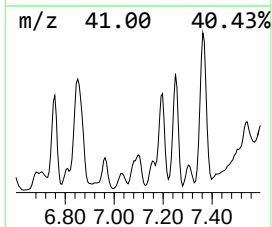
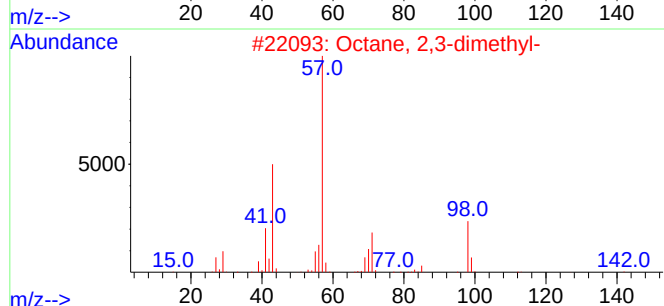
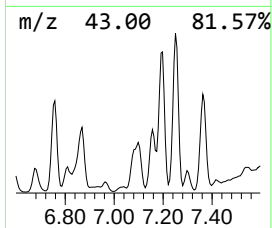
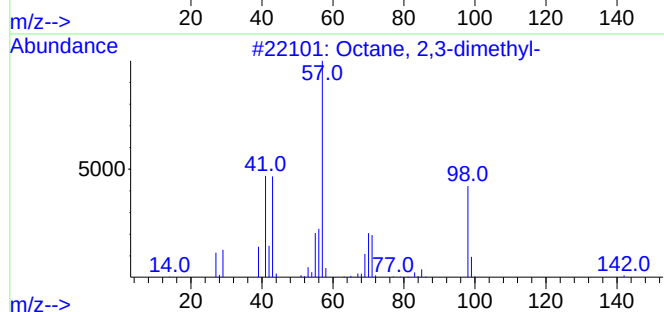
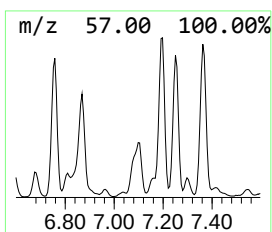
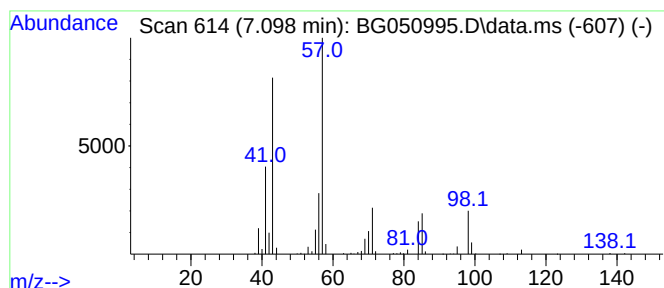
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.098	15.16 ng/ul	819640	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,3-dimethyl-			142	C10H22	007146-60-3	59
2	Octane, 2,3-dimethyl-			142	C10H22	007146-60-3	58
3	Heptane, 3-ethyl-2-methyl-			142	C10H22	014676-29-0	52
4	Nonane			128	C9H20	000111-84-2	50
5	Hexane, 2,4-dimethyl-			114	C8H18	000589-43-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

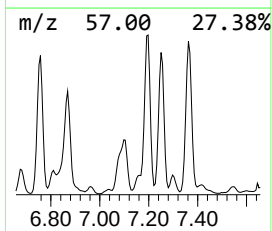
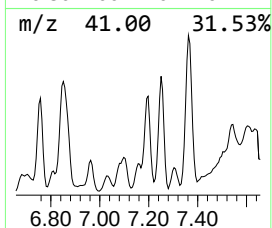
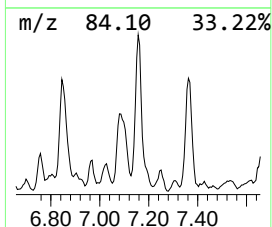
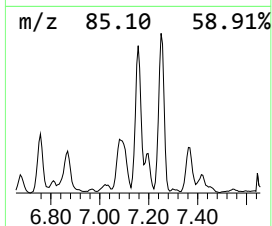
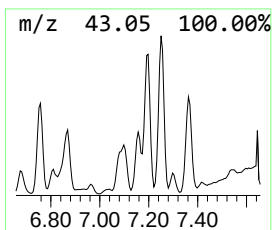
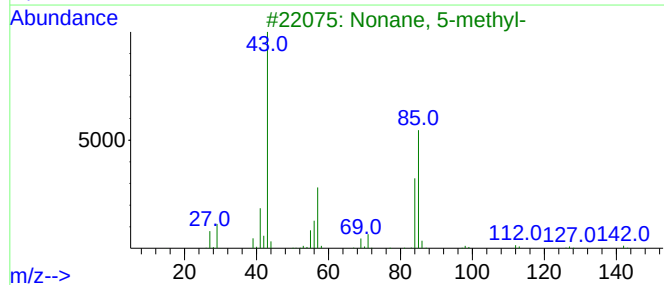
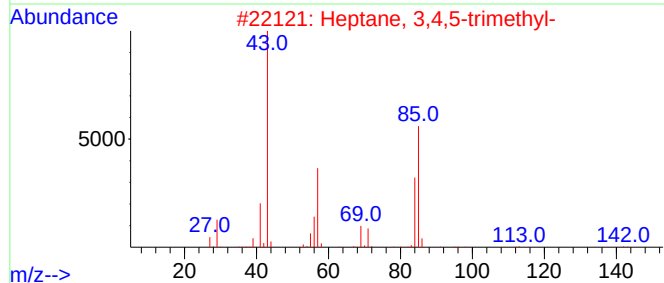
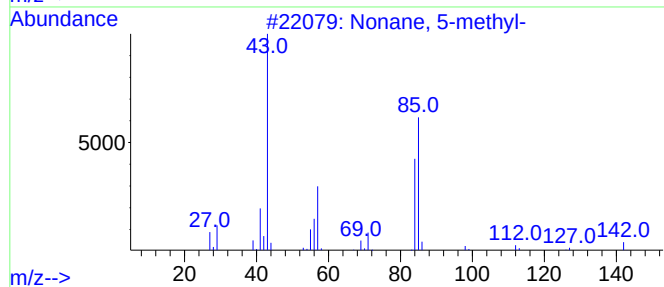
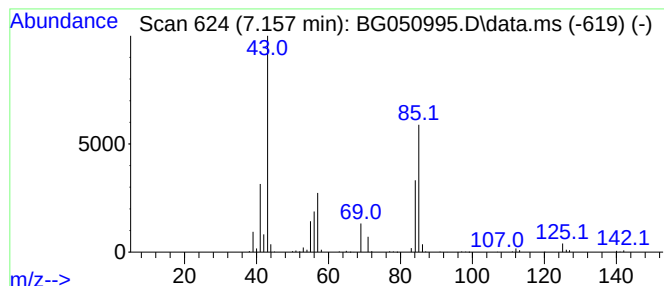
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.157	8.59 ng/ul	464183	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 5-methyl-	142	C10H22	015869-85-9	90
2		Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	86
3		Nonane, 5-methyl-	142	C10H22	015869-85-9	78
4		Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	72
5		Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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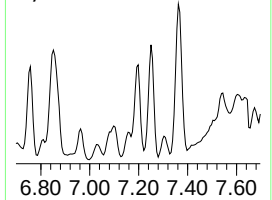
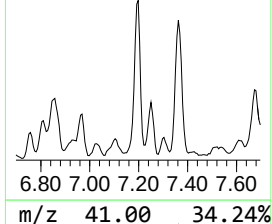
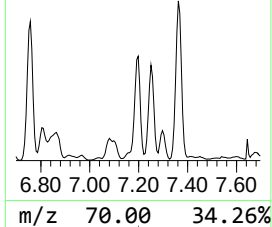
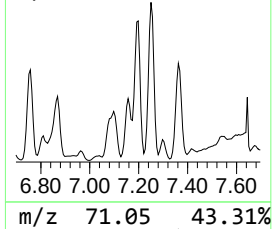
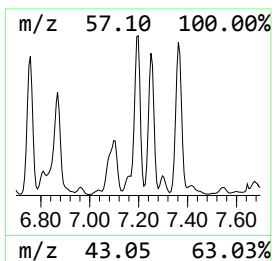
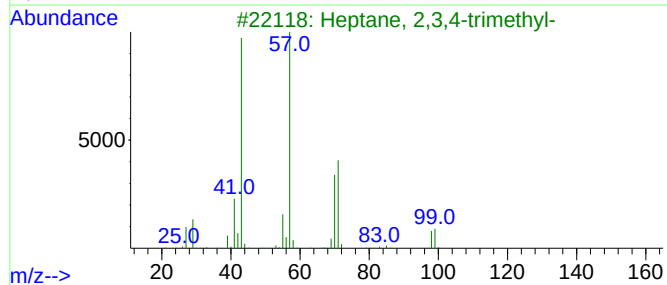
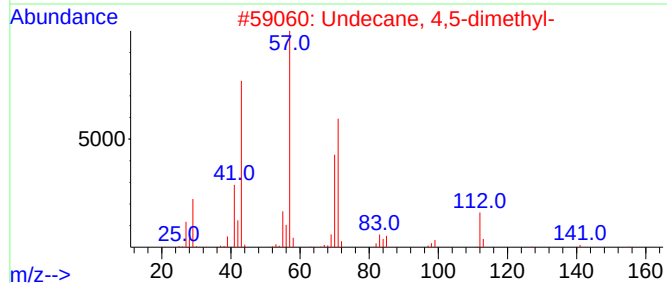
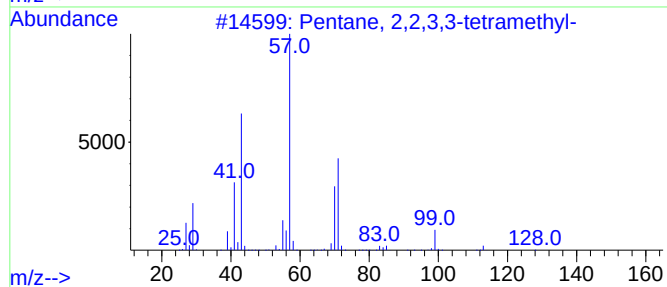
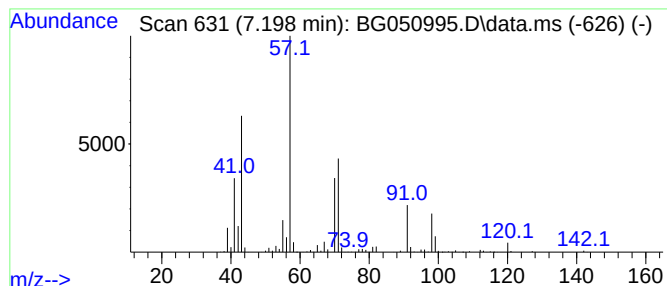
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.198	35.27 ng/ul	1906650	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	59
2		Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	59
3		Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	59
4		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	50
5		Nonane, 4-methyl-	142	C10H22	017301-94-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

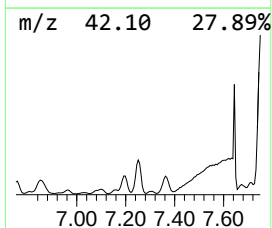
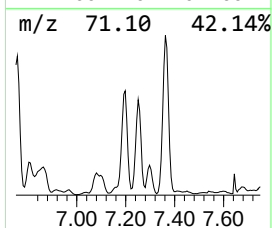
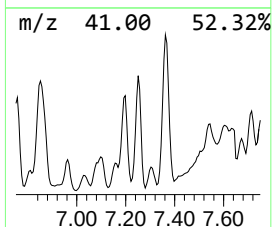
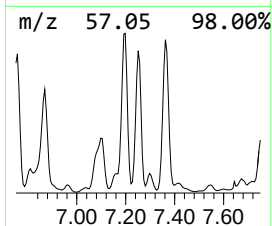
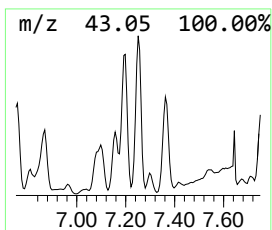
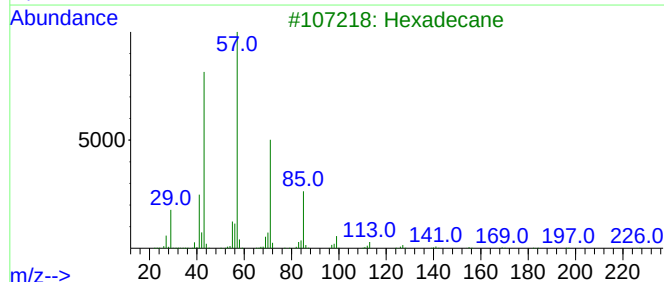
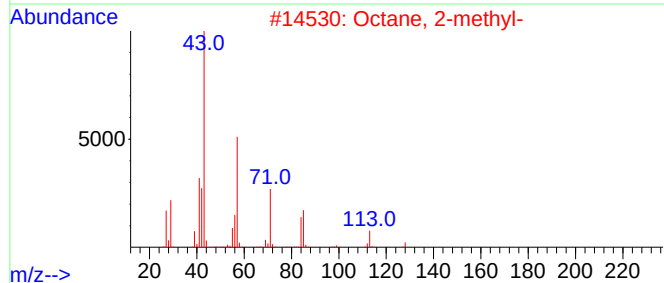
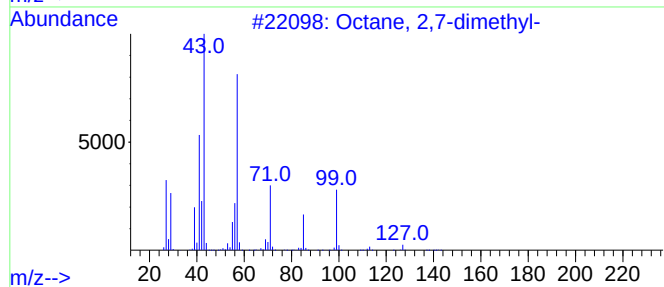
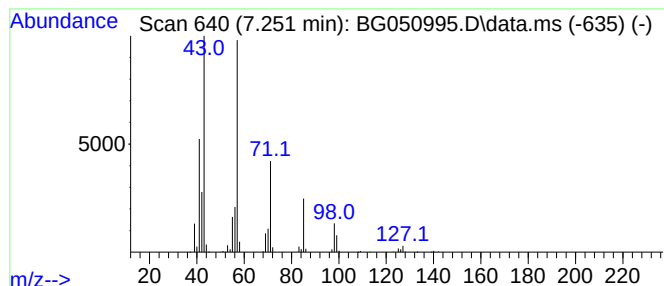
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.251	30.43 ng/ul	1645190	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	72
2		Octane, 2-methyl-	128	C9H20	003221-61-2	72
3		Hexadecane	226	C16H34	000544-76-3	64
4		Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	64
5		Tridecane, 6-methyl-	198	C14H30	013287-21-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

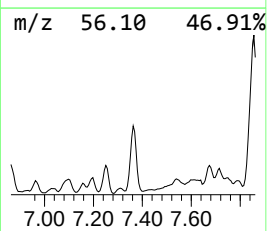
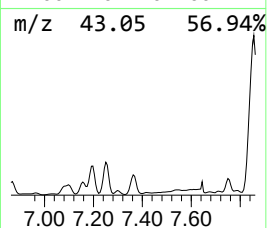
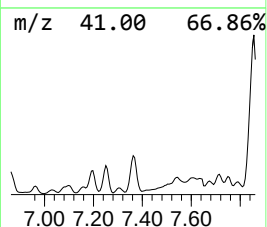
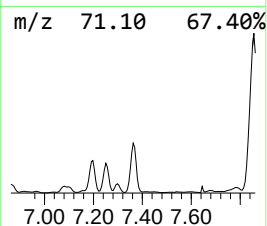
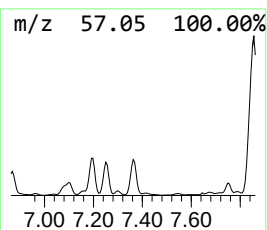
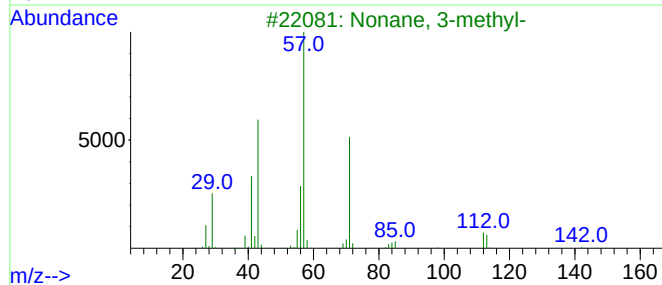
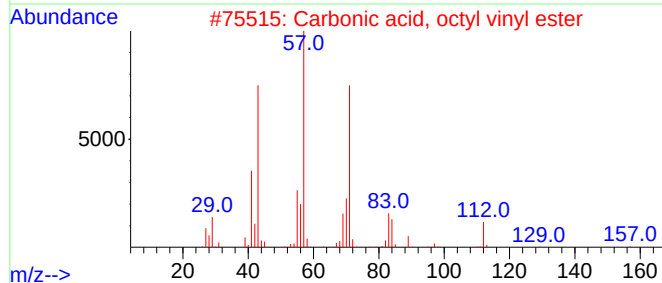
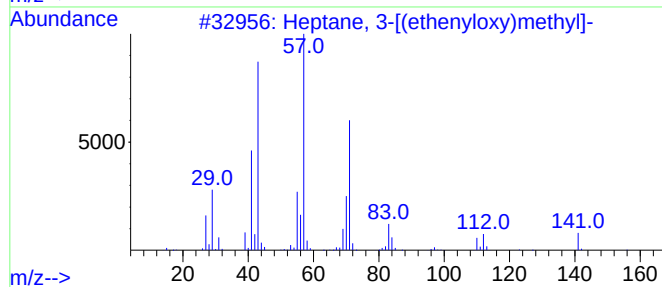
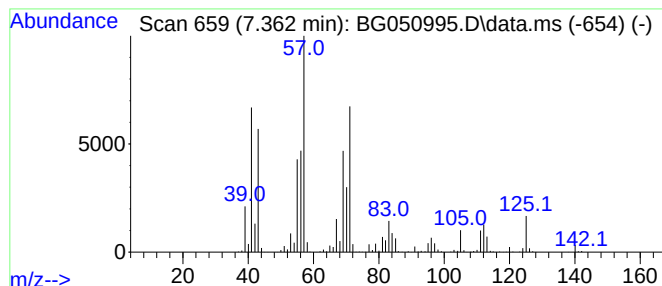
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 unknown-01 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.362	61.11 ng/ul	3303380	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-[(ethenyoxy)methyl]-	156	C10H20O	000103-44-6	47
2			Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	43
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	43
4			1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	43
5			Octane, 3-ethyl-	142	C10H22	005881-17-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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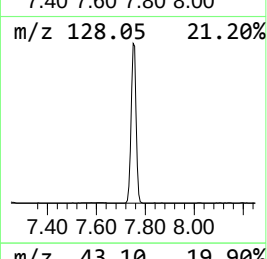
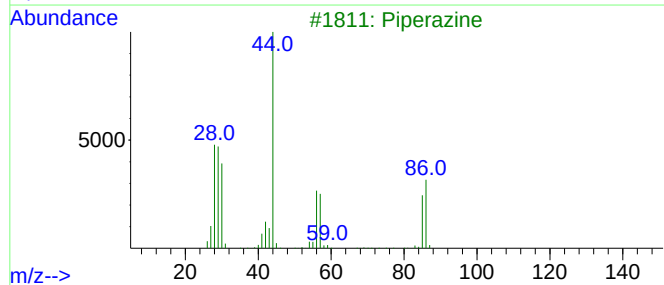
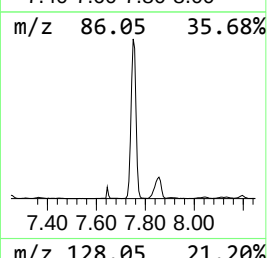
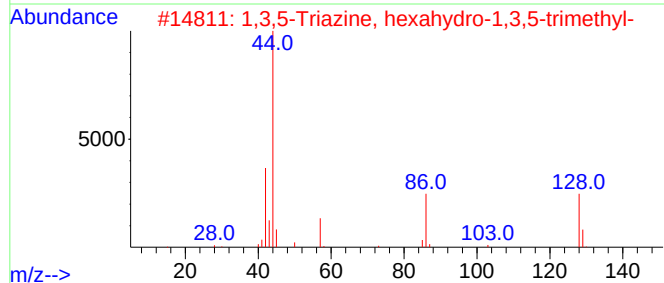
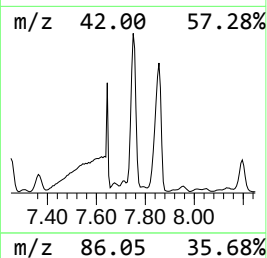
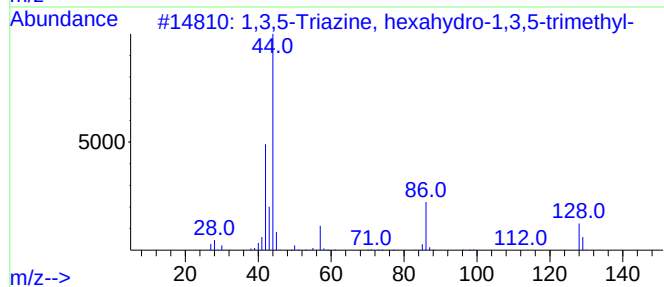
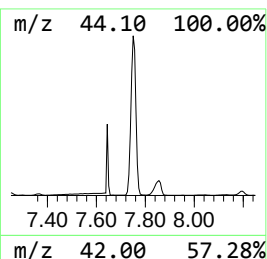
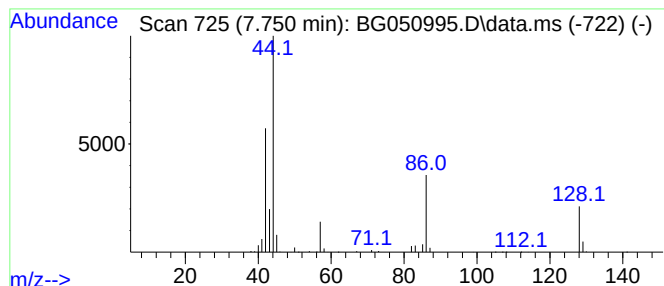
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 1,3,5-Triazine, hexahydro-1... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.750	51.06 ng/ul	2760500	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine, hexahydro-1,3,5-...	129	C6H15N3	000108-74-7	86		
2	1,3,5-Triazine, hexahydro-1,3,5-...	129	C6H15N3	000108-74-7	47		
3	Piperazine	86	C4H10N2	000110-85-0	9		
4	Piperazine	86	C4H10N2	000110-85-0	9		
5	Barbituric acid	128	C4H4N2O3	000067-52-7	7		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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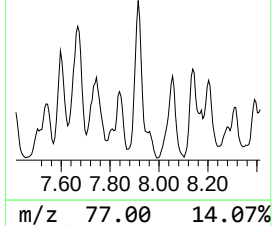
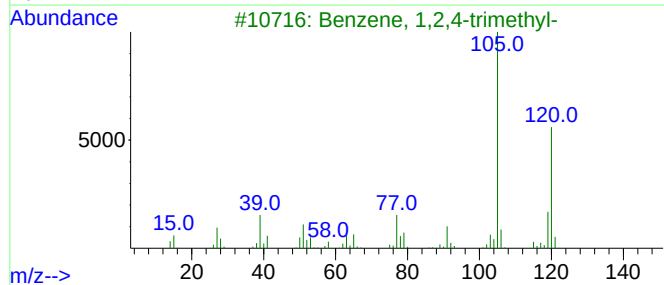
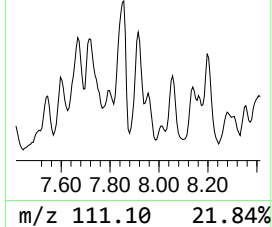
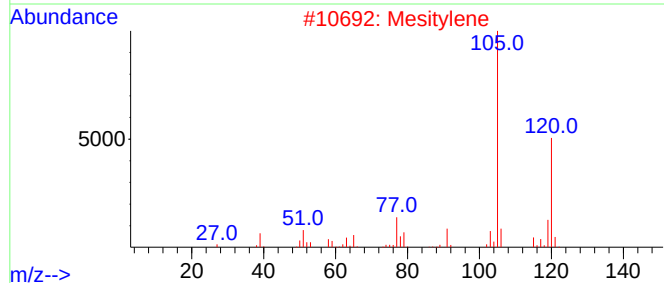
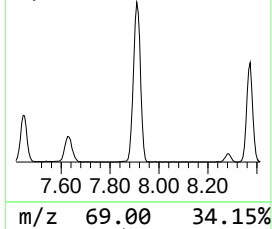
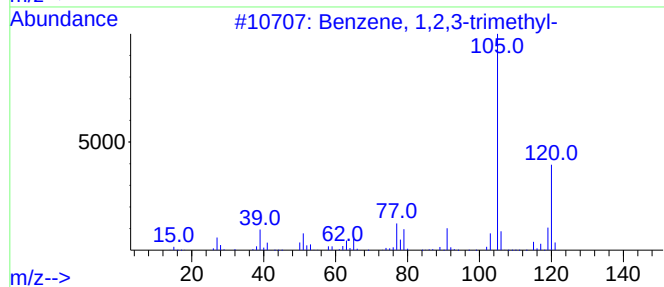
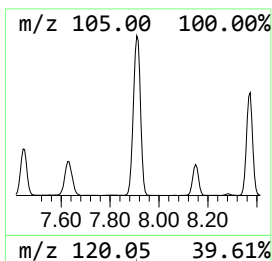
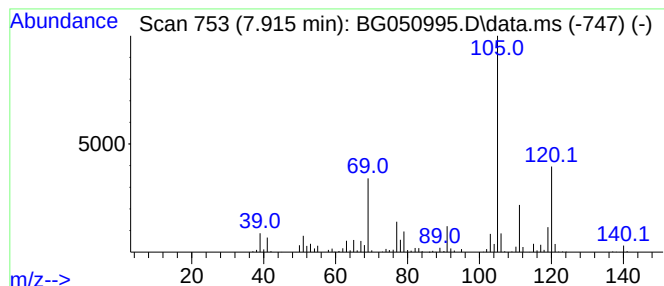
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Benzene, 1,2,3-trimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.915	33.25 ng/ul	1797300	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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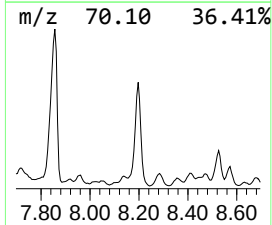
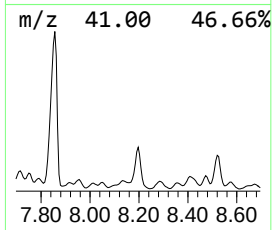
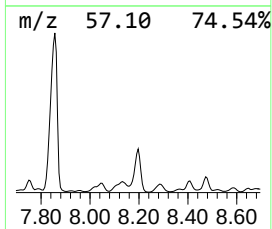
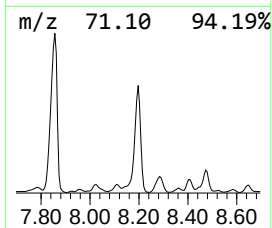
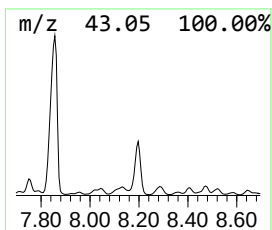
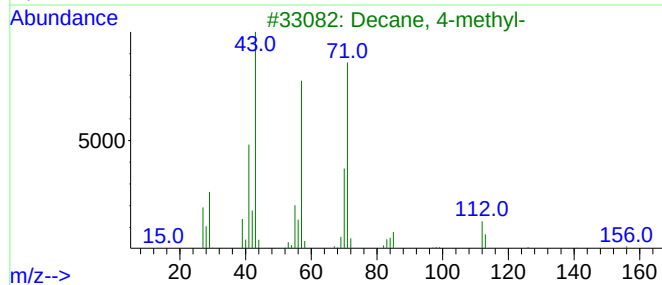
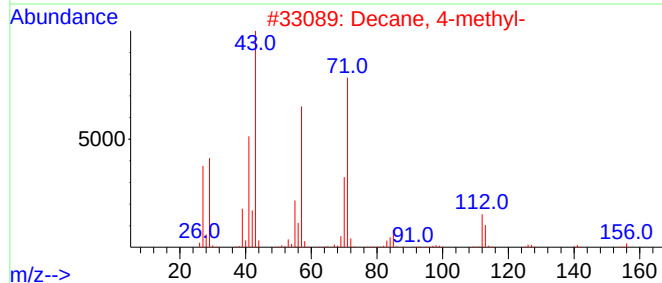
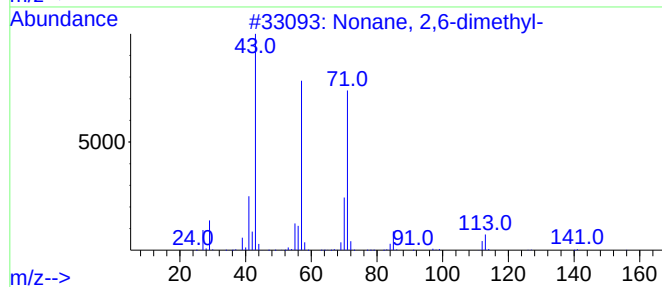
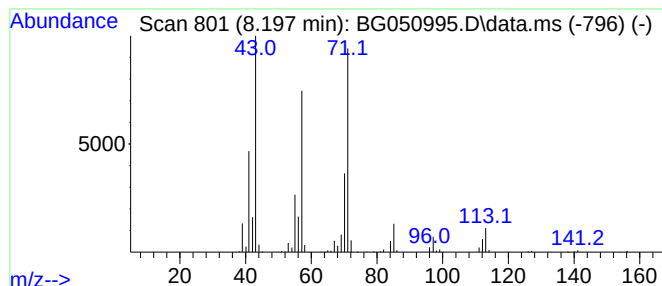
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.197	64.47 ng/ul	3485300	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	91
2		Decane, 4-methyl-	156	C11H24	002847-72-5	83
3		Decane, 4-methyl-	156	C11H24	002847-72-5	83
4		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
5		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

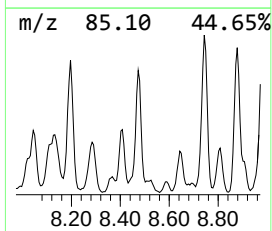
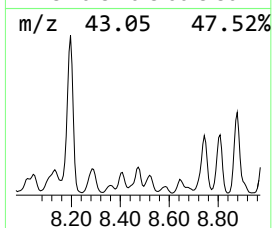
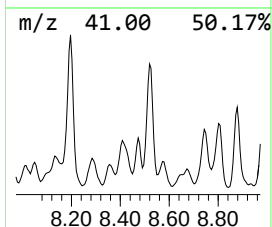
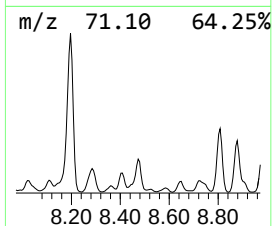
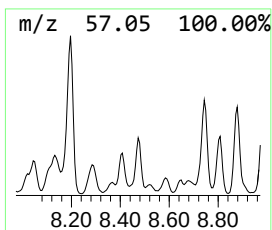
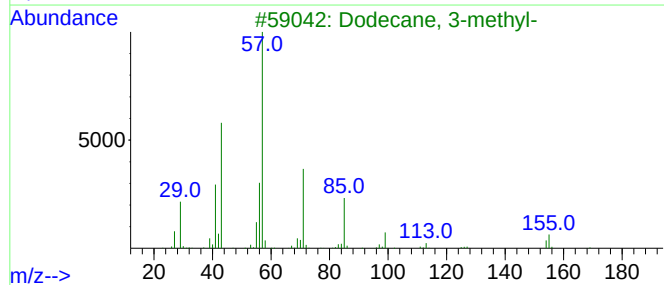
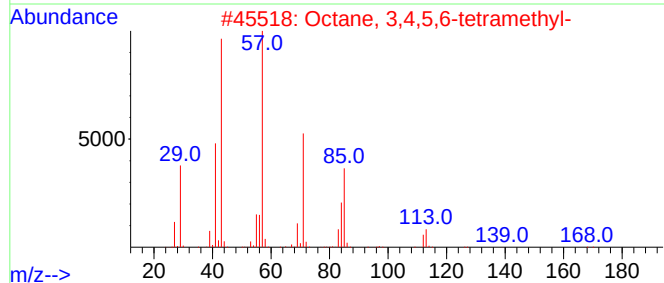
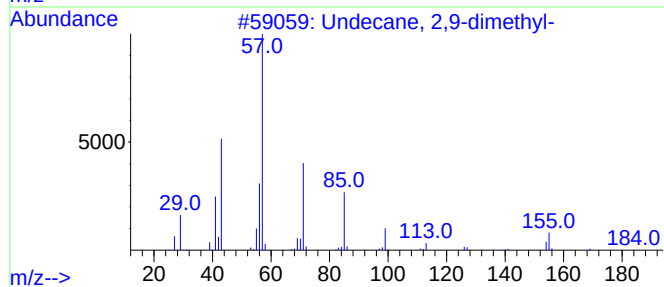
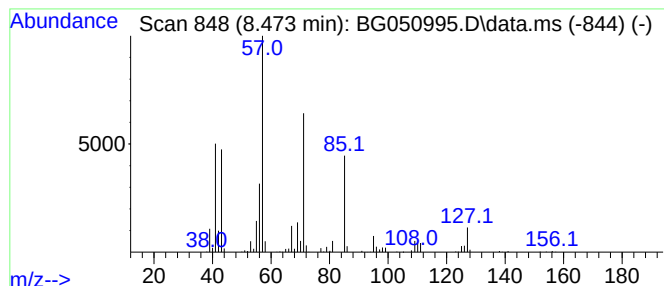
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.473	16.73 ng/ul	904580	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	64
2		Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	59
3		Dodecane, 3-methyl-	184	C13H28	017312-57-1	59
4		Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	53
5		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

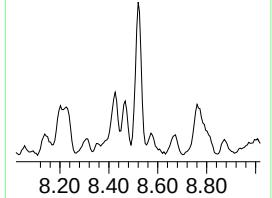
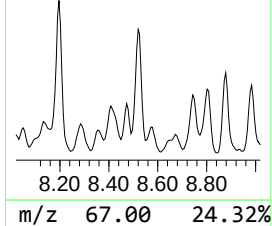
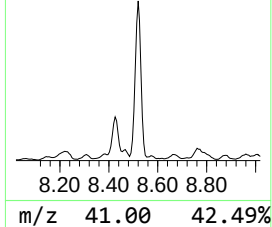
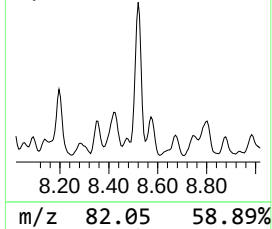
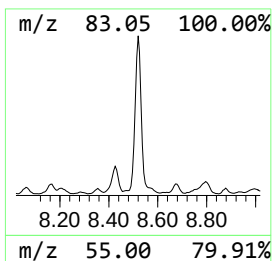
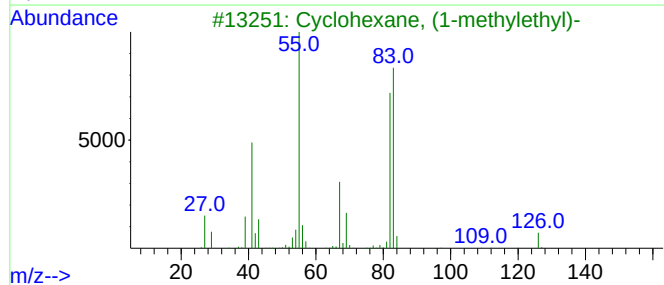
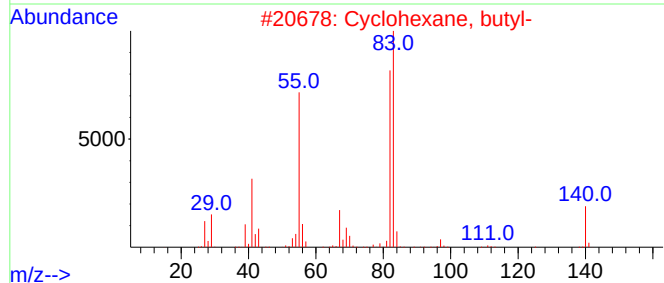
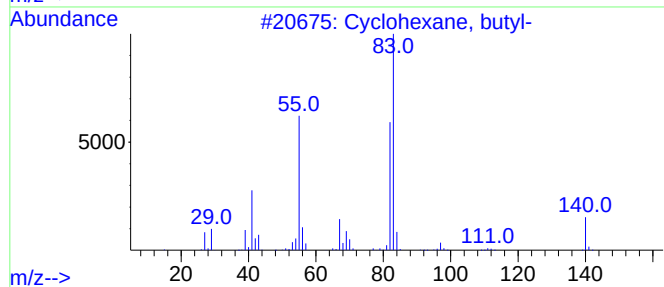
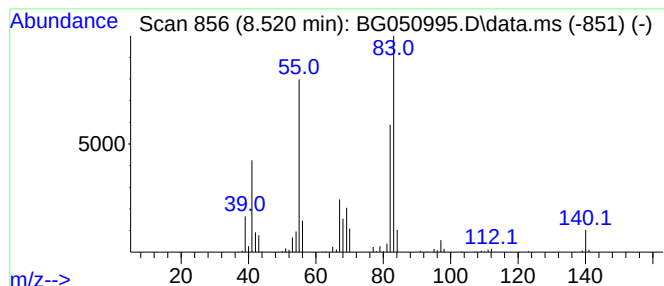
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 (DEL) Alkane: Cyclic8.520 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.520	49.83 ng/ul	2693930	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	90
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	74
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64
4			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	53
5			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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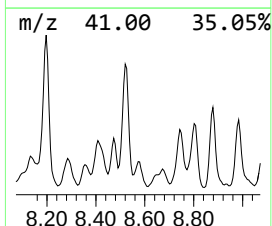
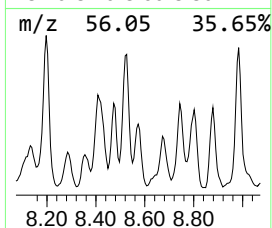
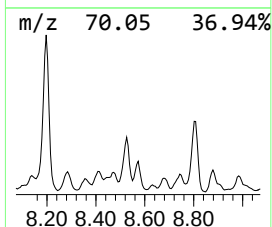
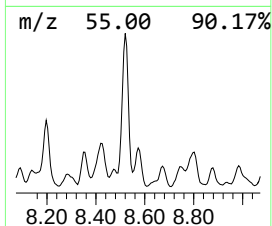
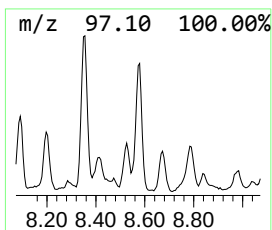
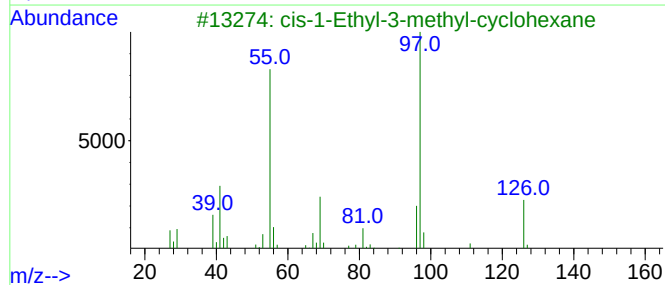
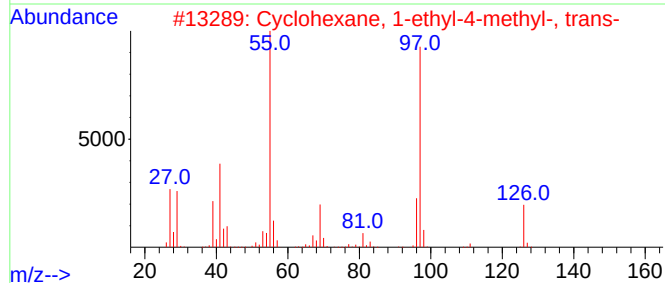
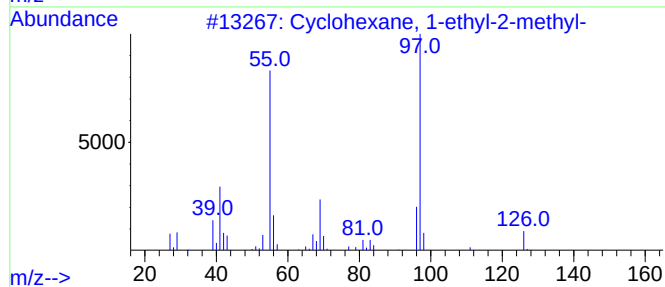
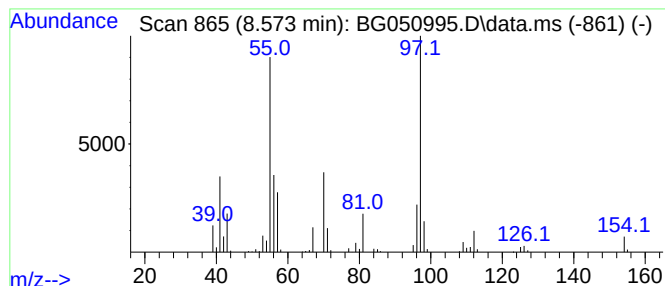
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 (DEL) Alkane: Cyclic8.573 Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.573	12.08 ng/ul	653201	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	59
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	59
3		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	53
4		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	53
5		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

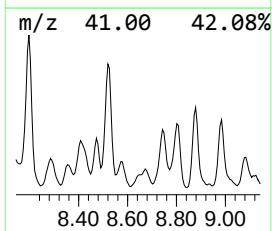
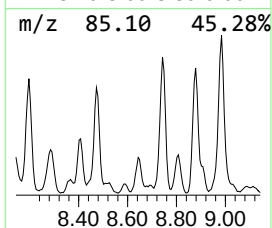
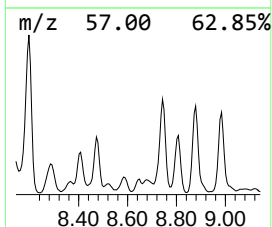
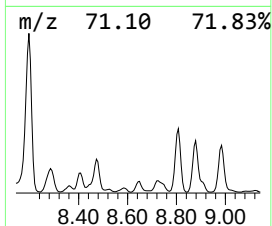
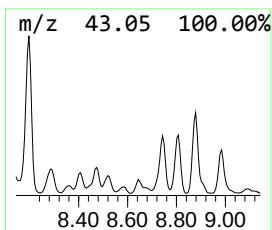
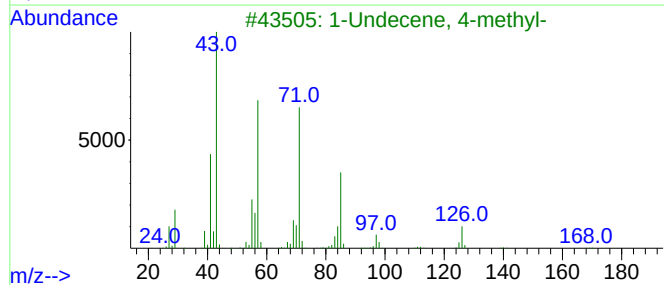
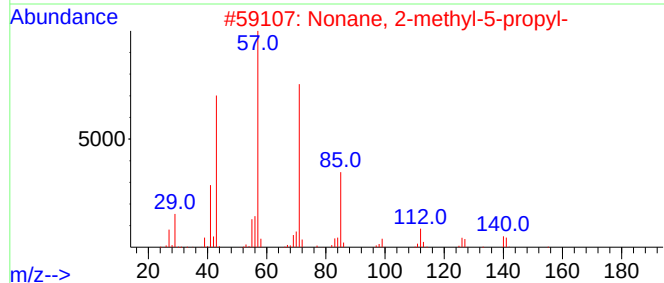
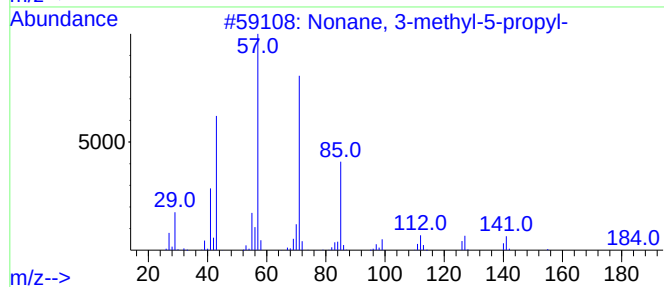
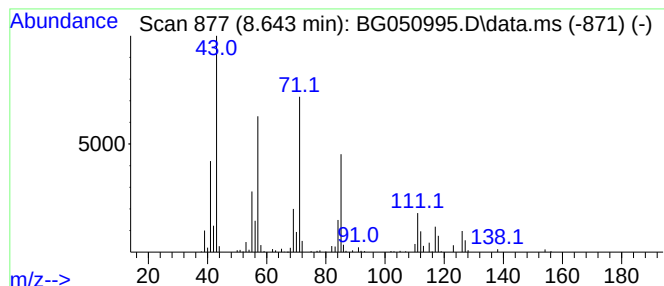
Instrument :
BNA_G
ClientSampleId :
COV04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 76

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.643	5.34 ng/ul	288837	1,4-Dichlorobenzene-d4			8.261
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-5-propyl-		184	C13H28	031081-18-2	59
2	Nonane, 2-methyl-5-propyl-		184	C13H28	031081-17-1	59
3	1-Undecene, 4-methyl-		168	C12H24	074630-39-0	53
4	Nonane, 5-butyl-		184	C13H28	017312-63-9	50
5	Decane, 4-ethyl-		170	C12H26	001636-44-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

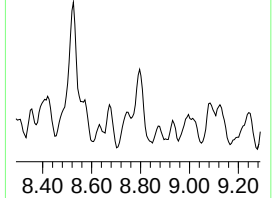
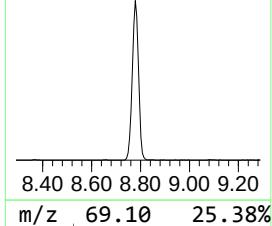
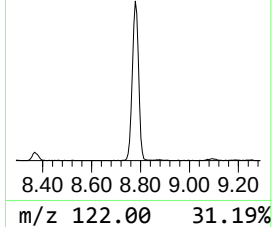
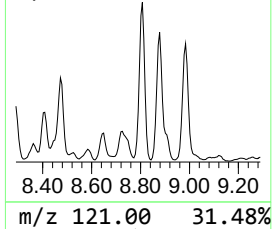
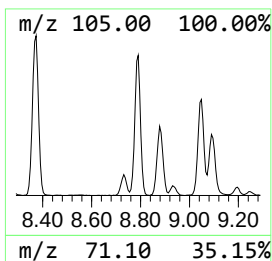
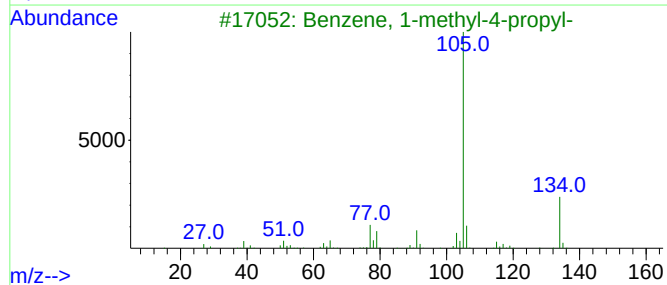
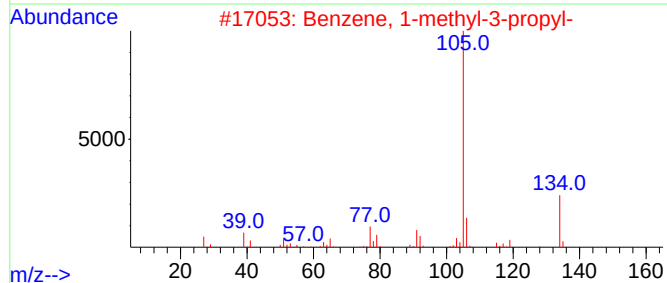
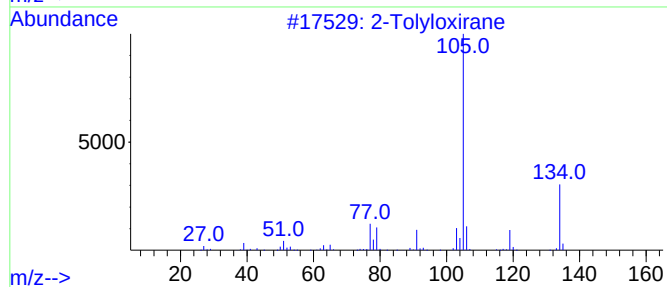
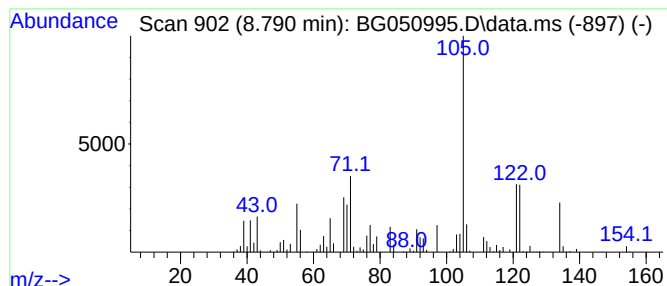
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 unknown-02 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.790	55.55 ng/ul	3002750	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Tolyloxirane		134	C9H10O	002783-26-8	38
2	Benzene, 1-methyl-3-propyl-		134	C10H14	001074-43-7	38
3	Benzene, 1-methyl-4-propyl-		134	C10H14	001074-55-1	38
4	Benzene, 1-methyl-2-propyl-		134	C10H14	001074-17-5	30
5	Benzene, 1-methyl-3-propyl-		134	C10H14	001074-43-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

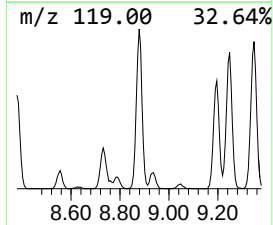
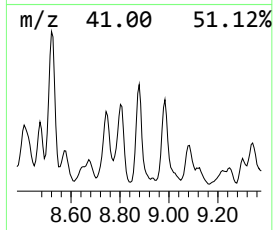
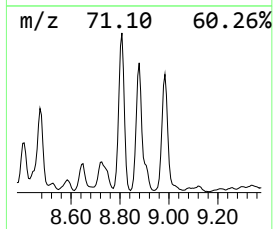
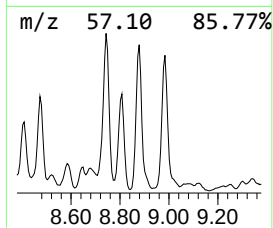
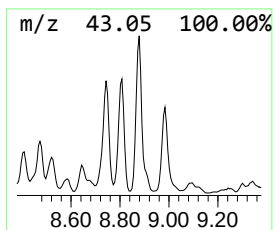
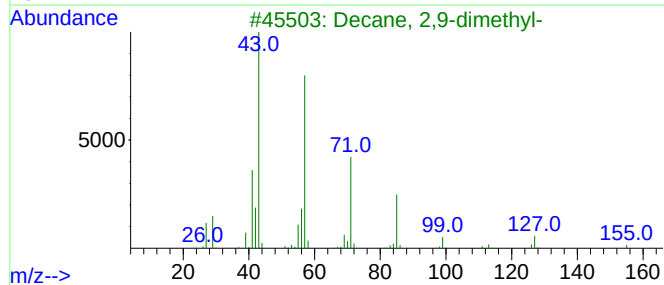
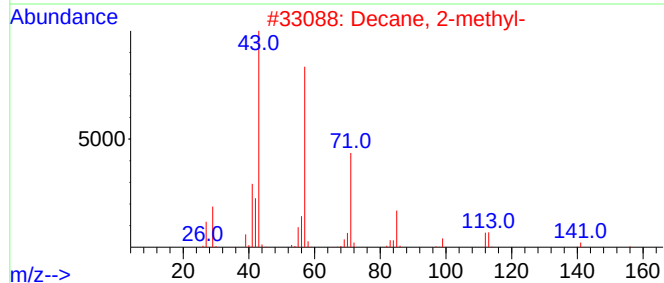
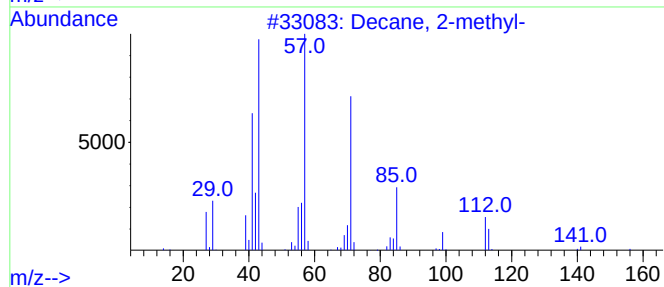
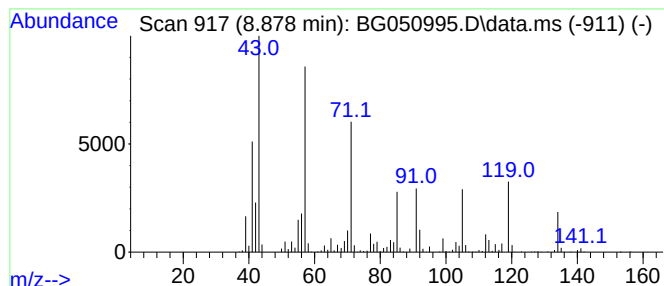
Instrument :
BNA_G
ClientSampleId :
COV04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.878	39.23 ng/ul	2120690	1,4-Dichlorobenzene-d4	8.261	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2-methyl-	156	C11H24	006975-98-0	49
2	Decane, 2-methyl-	156	C11H24	006975-98-0	46
3	Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	43
4	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	41
5	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

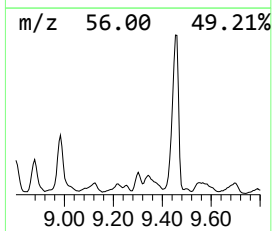
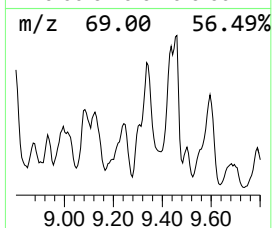
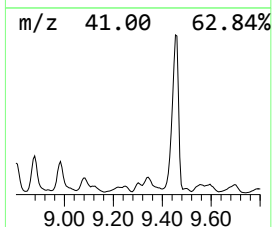
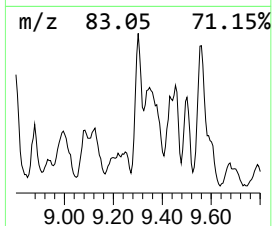
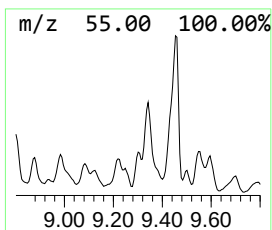
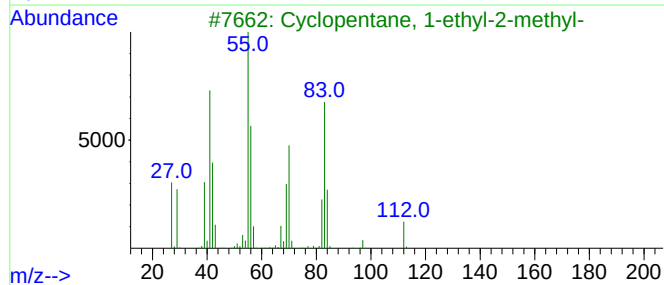
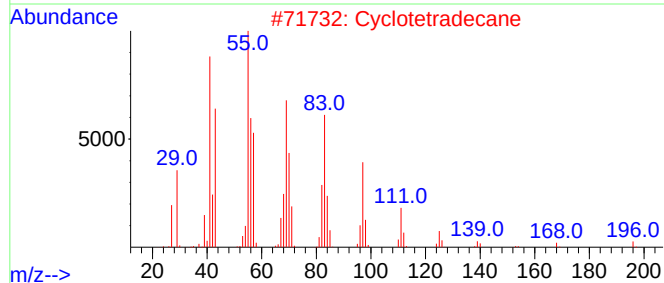
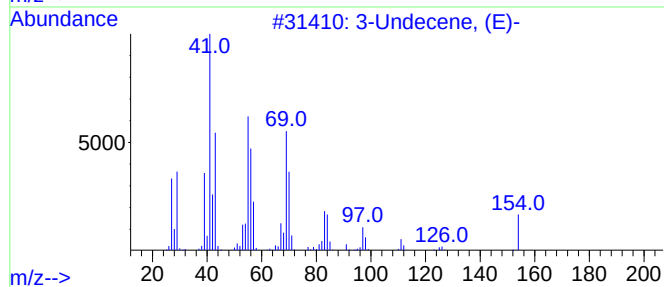
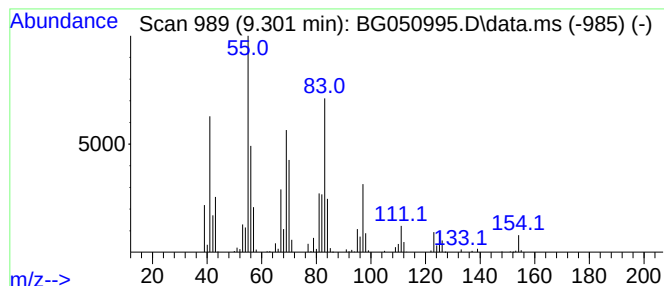
Instrument :
BNA_G
ClientSampleId :
COV04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 72

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.301	7.82 ng/ul	422680	1,4-Dichlorobenzene-d4			8.261
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Undecene, (E)-		154	C11H22	001002-68-2	64
2	Cyclotetradecane		196	C14H28	000295-17-0	64
3	Cyclopentane, 1-ethyl-2-methyl-		112	C8H16	003726-46-3	62
4	4-Heptafluorobutyryloxyhexadecane		438	C20H33F7O2	1010282-97-2	62
5	3-Undecene, (Z)-		154	C11H22	000821-97-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

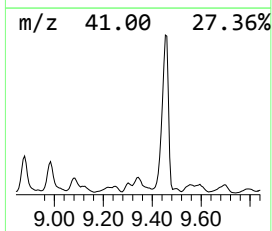
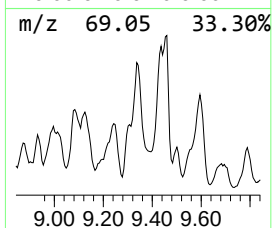
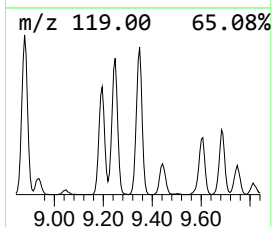
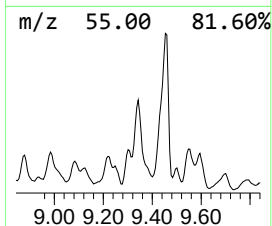
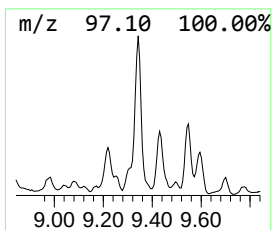
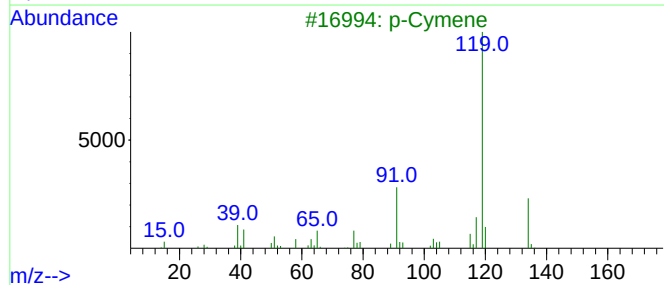
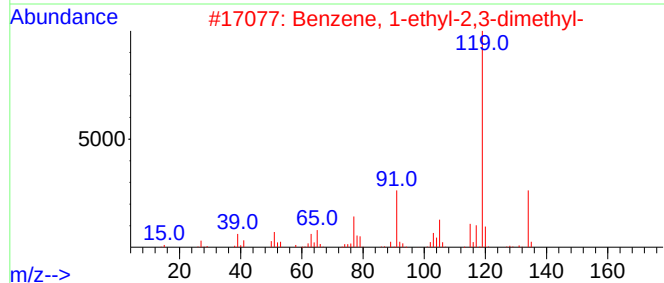
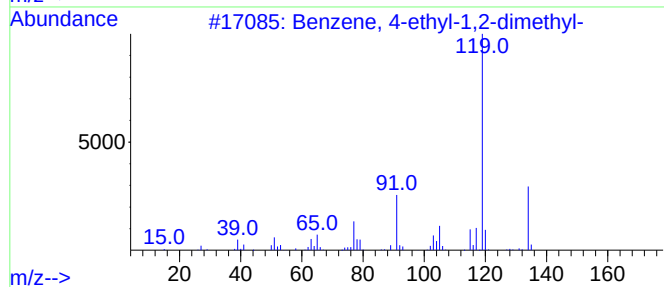
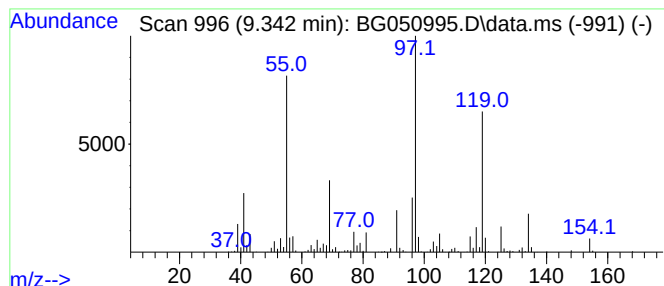
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.342	29.76 ng/ul	1608920	1,4-Dichlorobenzene-d4	8.261

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	78
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	78
3			p-Cymene	134	C10H14	000099-87-6	78
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	78
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

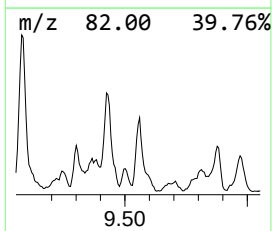
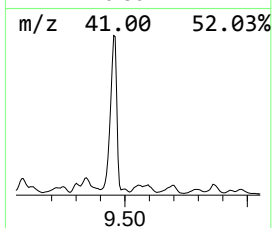
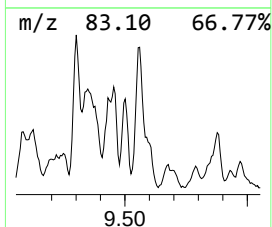
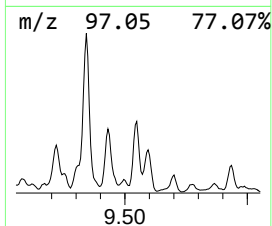
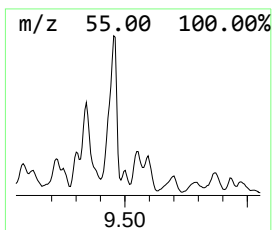
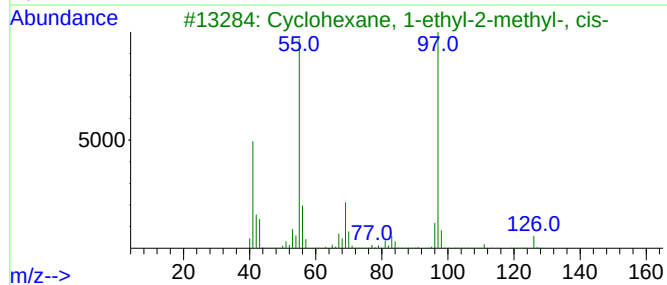
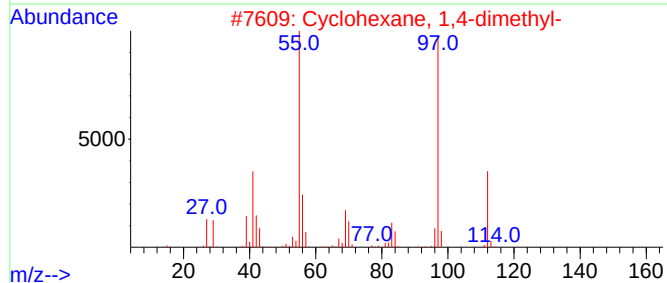
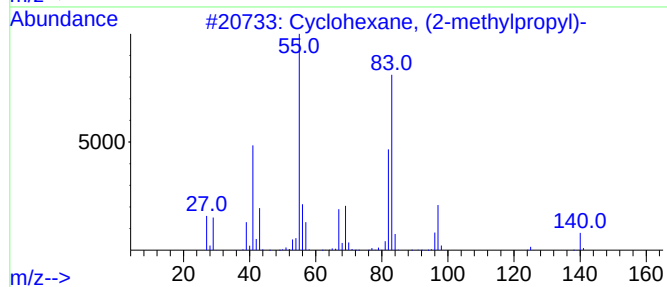
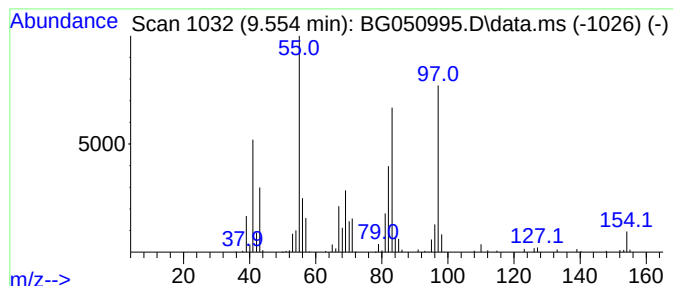
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 (DEL) Alkane: Cyclic9.554 Concentration Rank 71

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.554	8.01 ng/ul	433099	1,4-Dichlorobenzene-d4	8.261

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	50
2		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	47
3		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	46
4		Oxalic acid, cyclohexylmethyl et...	214	C11H18O4	1000309-68-0	43
5		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

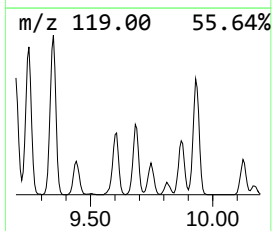
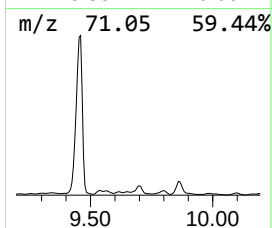
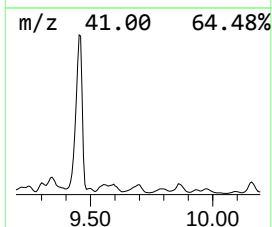
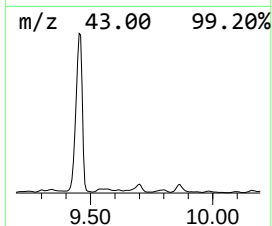
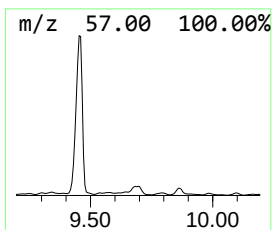
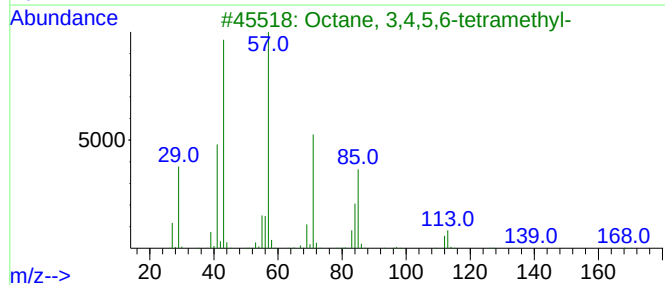
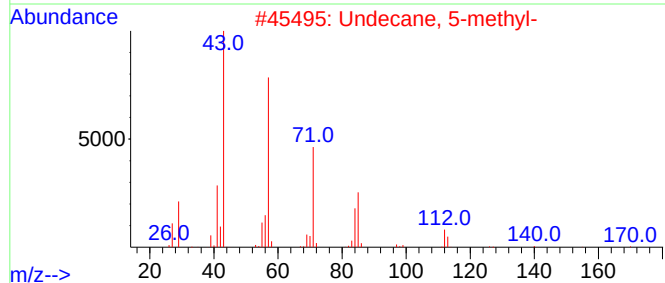
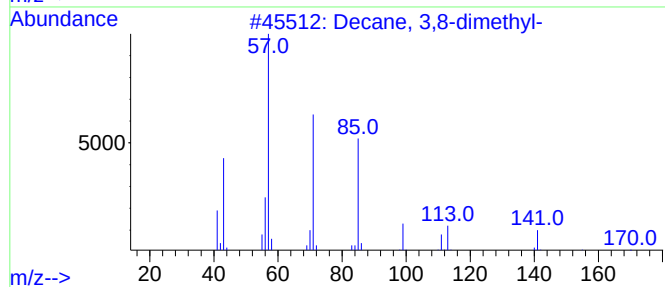
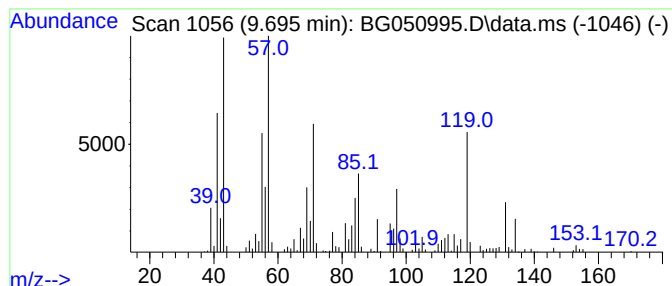
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.695	40.02 ng/ul	661174	Naphthalene-d8	11.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	42
2			Undecane, 5-methyl-	170	C12H26	001632-70-8	41
3			Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	38
4			Undecane, 5-methyl-	170	C12H26	001632-70-8	38
5			Octane, 2-methyl-	128	C9H20	003221-61-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

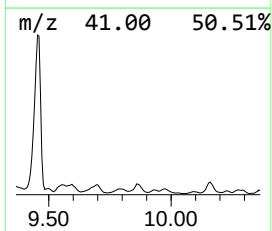
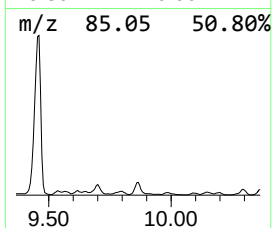
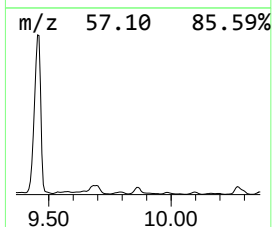
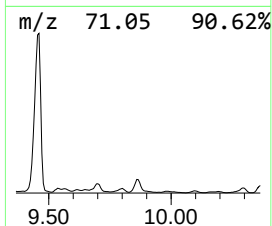
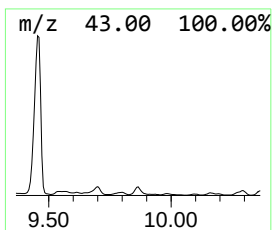
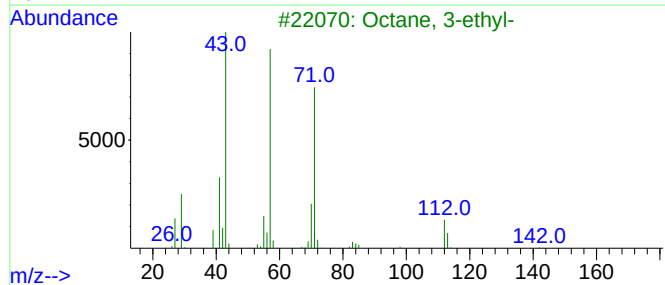
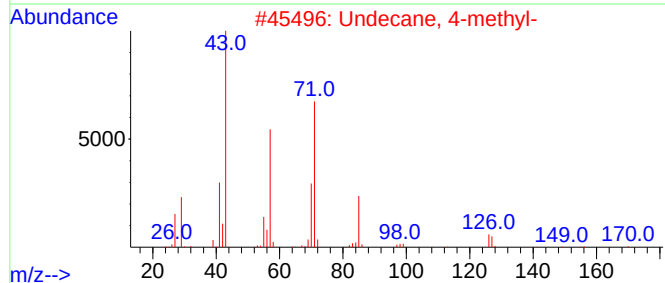
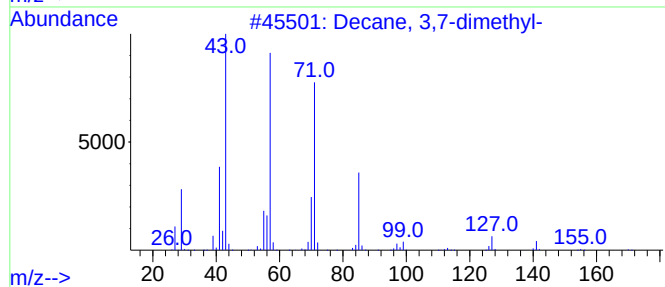
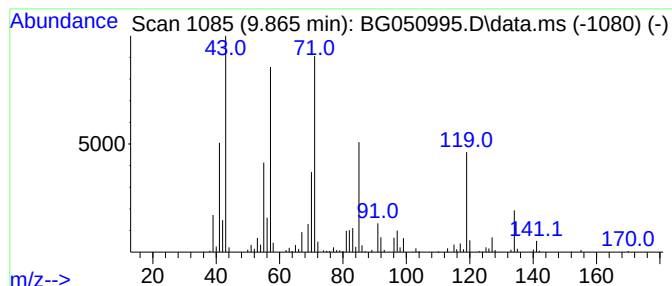
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.865	35.98 ng/ul	594514	Naphthalene-d8	11.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	60
2			Undecane, 4-methyl-	170	C12H26	002980-69-0	46
3			Octane, 3-ethyl-	142	C10H22	005881-17-4	43
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	38
5			Decane, 3-bromo-	220	C10H21Br	030571-71-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

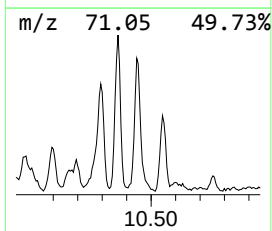
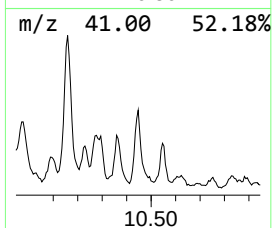
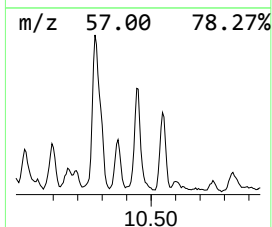
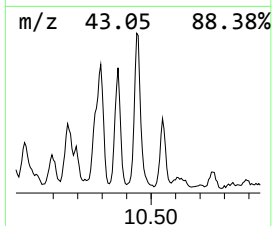
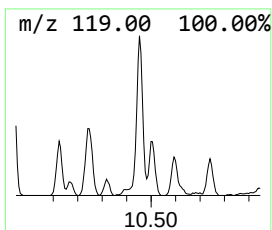
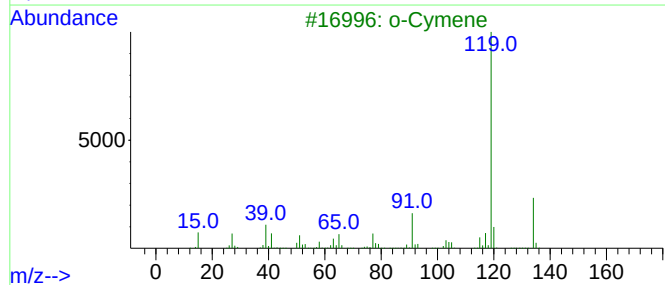
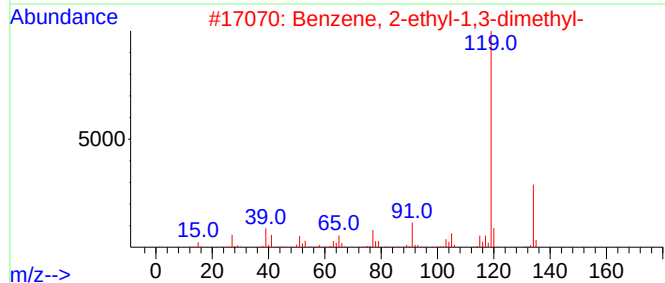
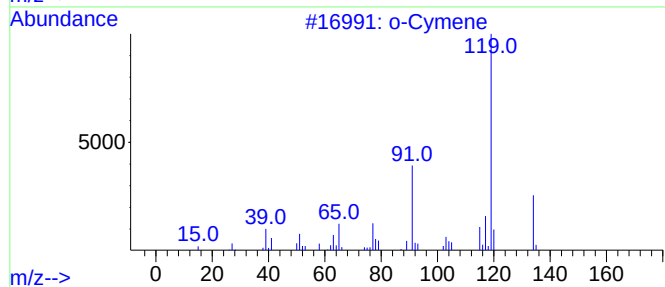
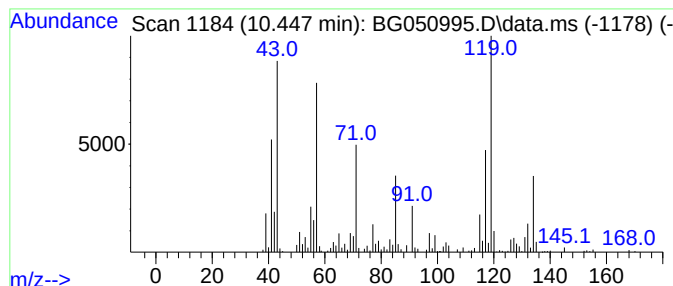
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 o-Cymene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.447	35.19 ng/ul	581320	Naphthalene-d8	11.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	50
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	46
3			o-Cymene	134	C10H14	000527-84-4	46
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	46
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
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Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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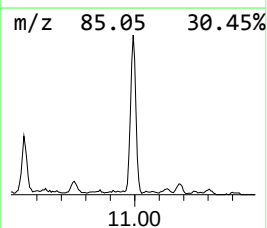
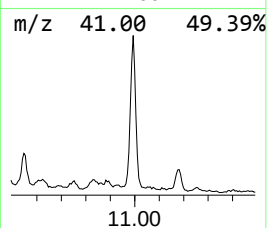
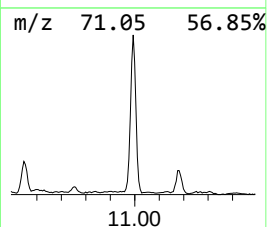
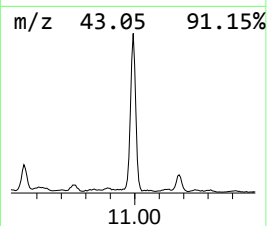
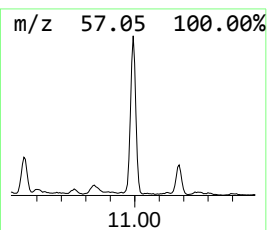
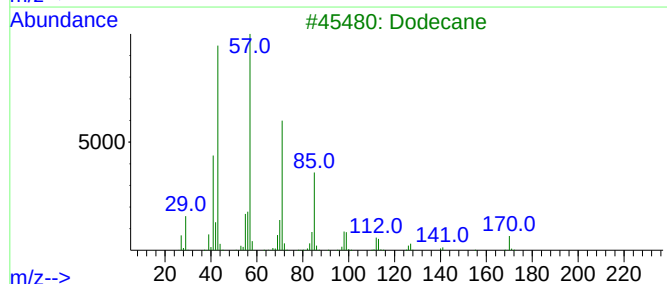
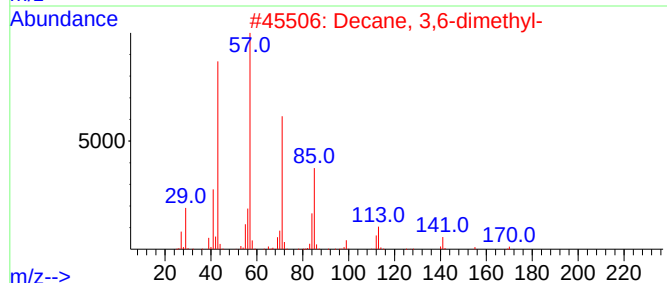
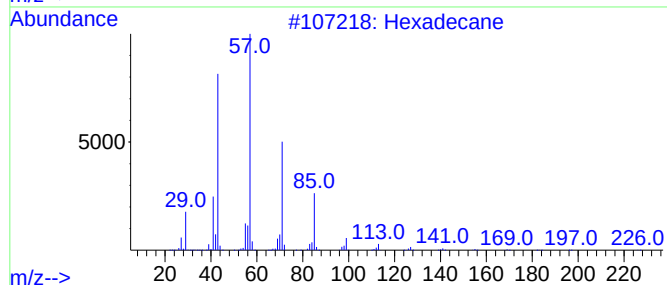
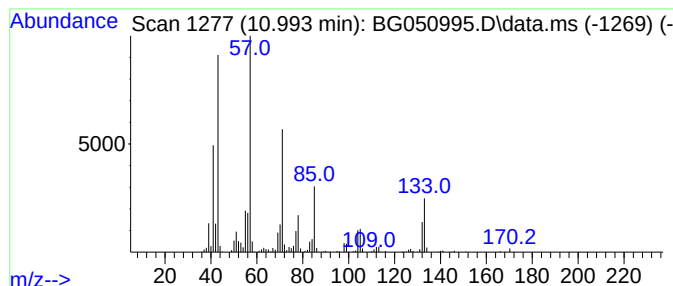
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.993	49.83 ng/ul	823322	Naphthalene-d8	11.064

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	49
2		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	49
3		Dodecane	170	C12H26	000112-40-3	49
4		Undecane, 4-methyl-	170	C12H26	002980-69-0	49
5		Tridecane	184	C13H28	000629-50-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

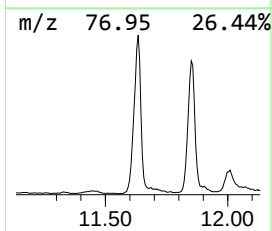
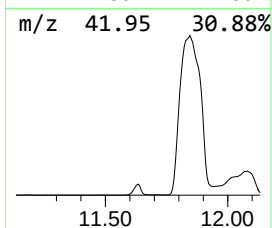
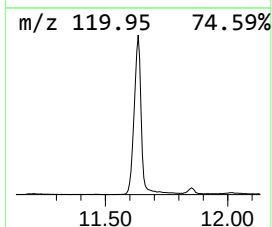
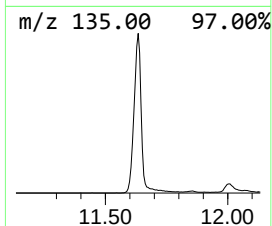
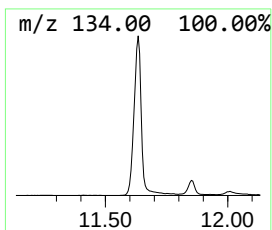
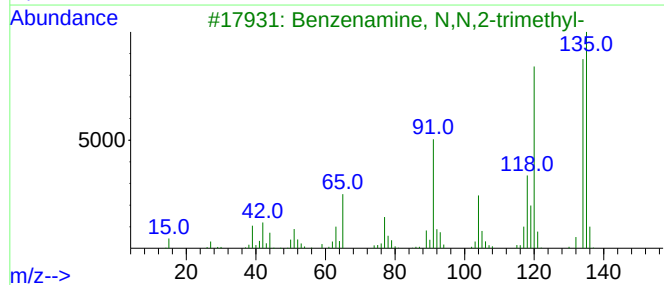
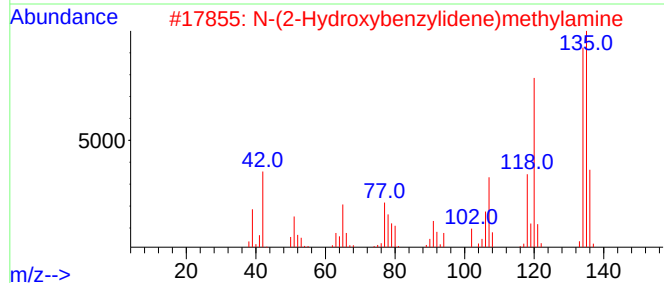
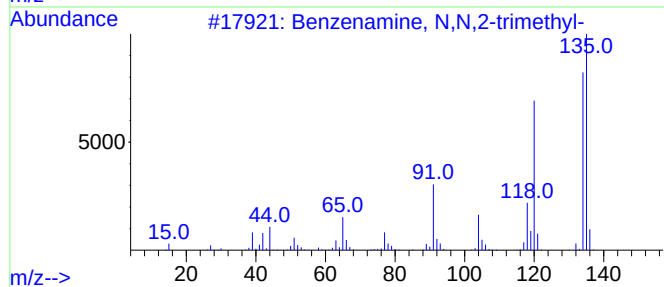
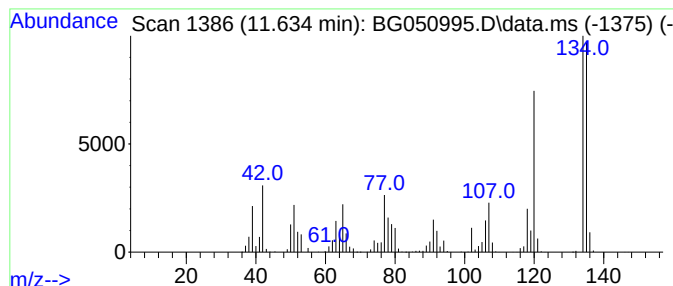
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 Benzenamine, N,N,2-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.634	231.58 ng/ul	3826050	Naphthalene-d8	11.064

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenamine, N,N,2-trimethyl-	135	C9H13N	000609-72-3	83
2		N-(2-Hydroxybenzylidene)methylamine	135	C8H9NO	1000433-53-5	81
3		Benzenamine, N,N,2-trimethyl-	135	C9H13N	000609-72-3	64
4		Benzenamine, N,N,3-trimethyl-	135	C9H13N	000121-72-2	53
5		6,7-Dihydro-5H-cyclopenta[d]pyri...	135	C7H9N3	026979-06-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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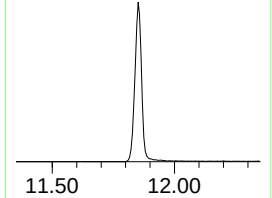
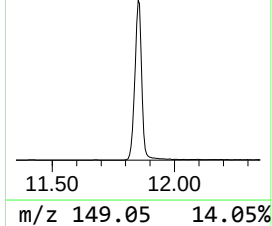
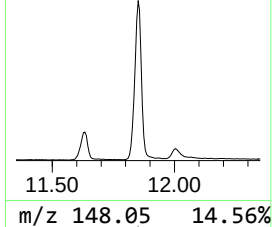
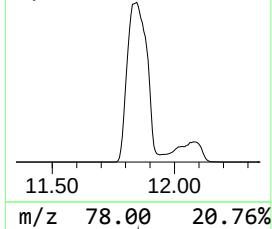
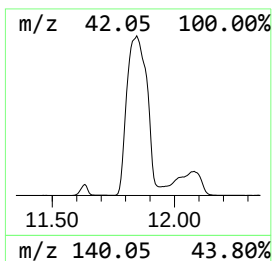
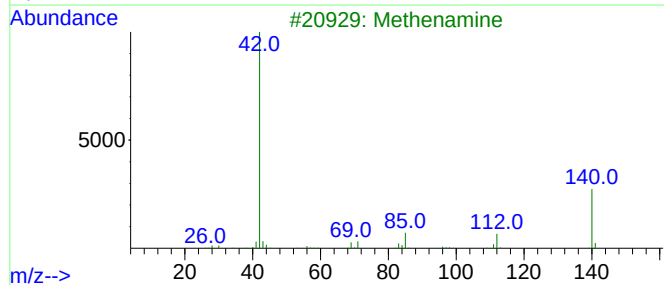
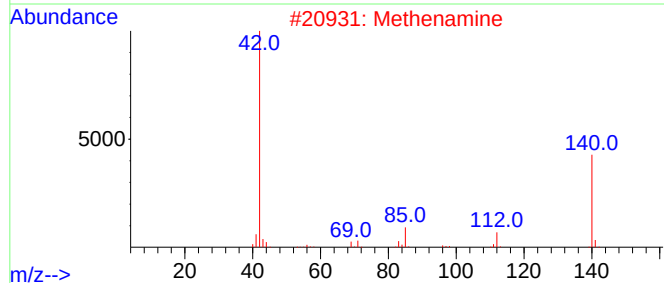
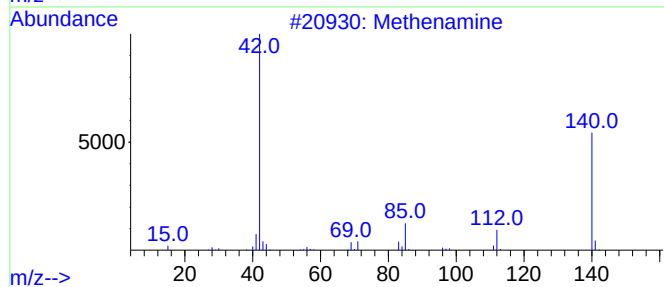
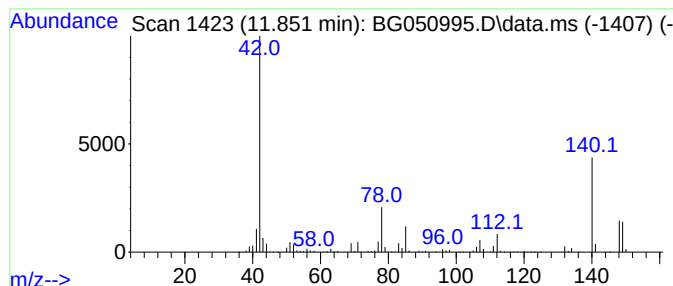
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 Methenamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.851	1001.70 ng/ul	16549600	Naphthalene-d8	11.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methenamine	140	C6H12N4	000100-97-0	94
2			Methenamine	140	C6H12N4	000100-97-0	93
3			Methenamine	140	C6H12N4	000100-97-0	76
4			Methenamine	140	C6H12N4	000100-97-0	76
5			Methenamine	140	C6H12N4	000100-97-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

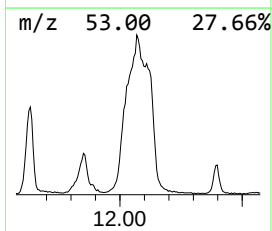
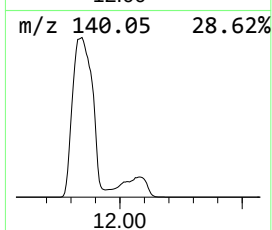
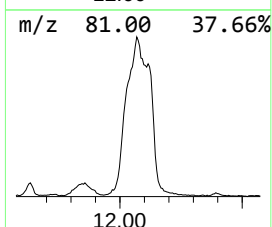
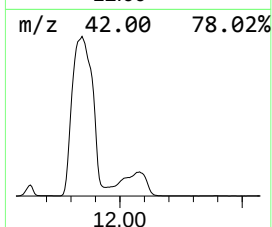
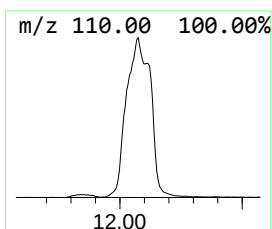
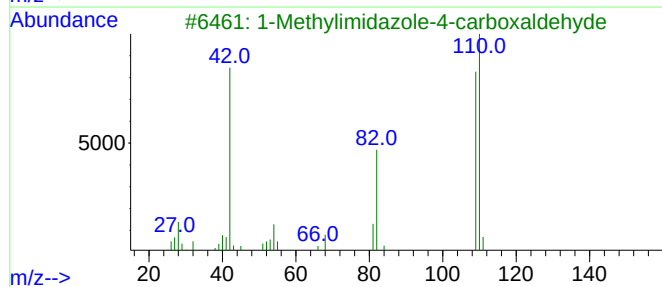
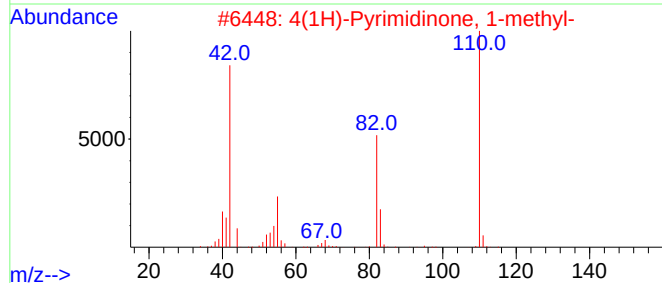
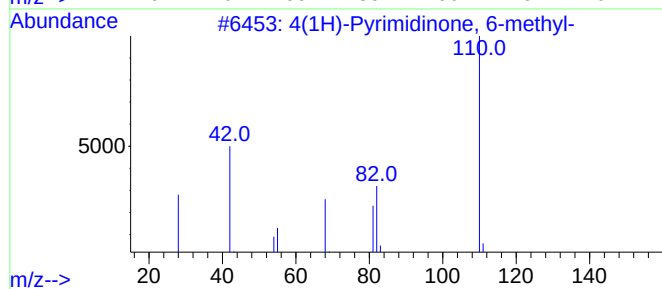
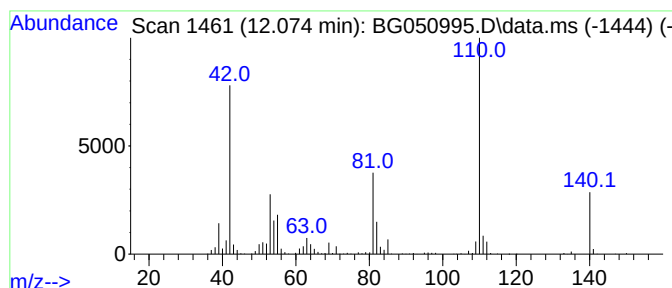
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 4(1H)-Pyrimidinone, 6-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.074	324.32 ng/ul	5358150	Naphthalene-d8	11.064

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4(1H)-Pyrimidinone, 6-methyl-	110	C5H6N2O	003524-87-6	59
2		4(1H)-Pyrimidinone, 1-methyl-	110	C5H6N2O	002228-30-0	58
3		1-Methylimidazole-4-carboxaldehyde	110	C5H6N2O	017289-26-8	53
4		Hydroquinone	110	C6H6O2	000123-31-9	50
5		Methenamine	140	C6H12N4	000100-97-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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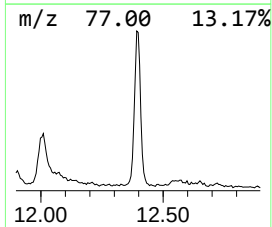
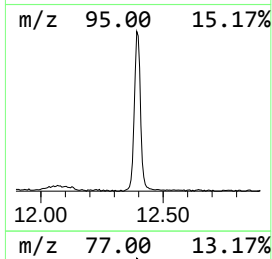
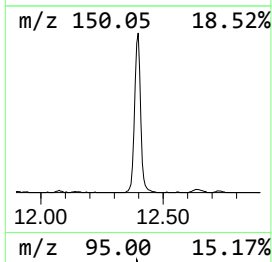
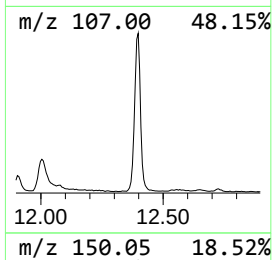
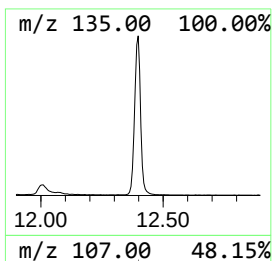
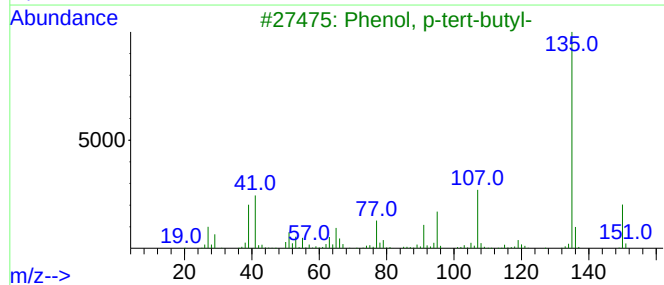
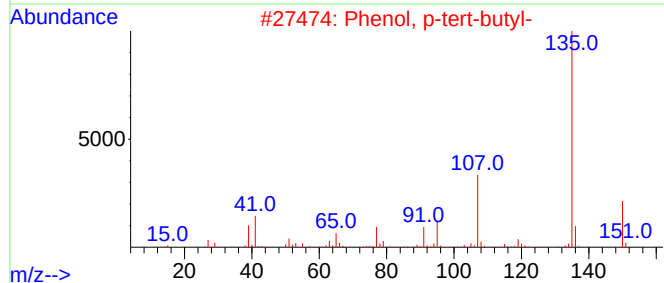
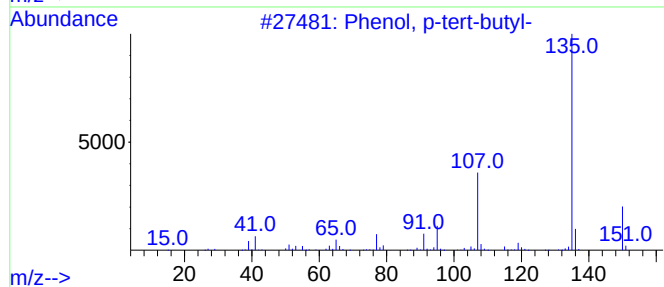
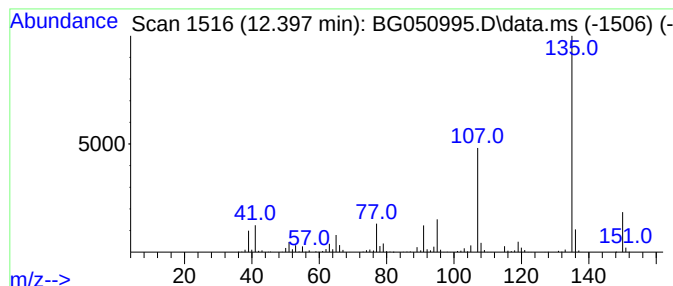
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.398	76.68 ng/ul	1266790	Naphthalene-d8	11.064

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	93
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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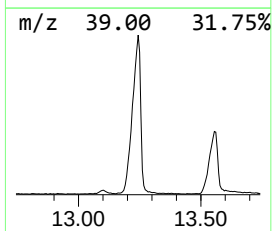
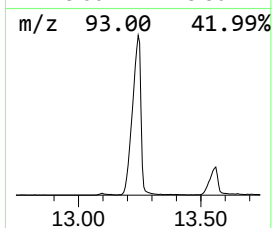
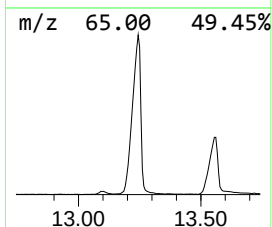
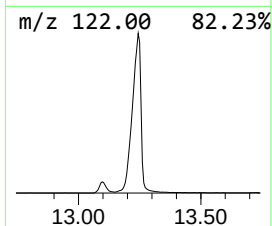
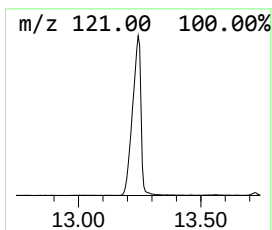
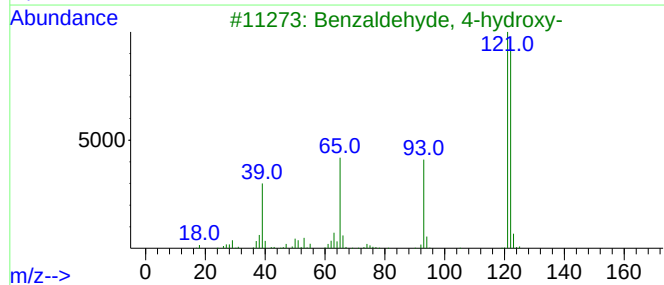
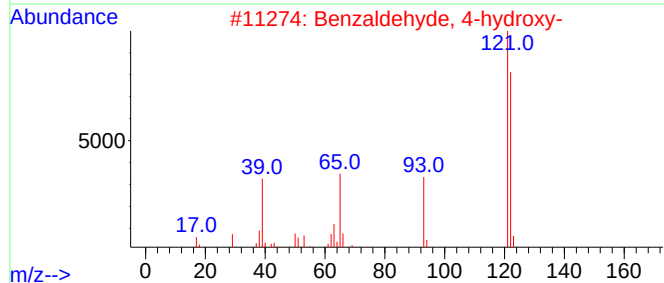
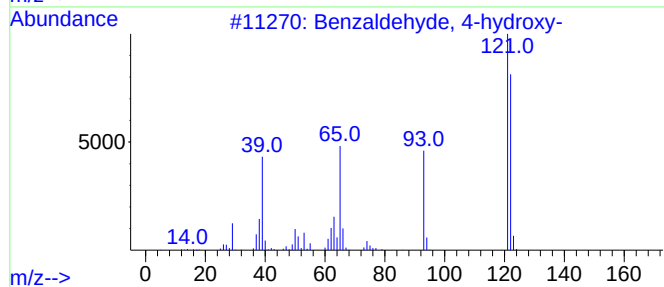
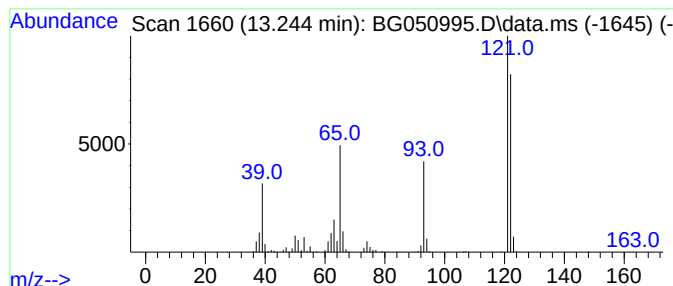
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 Benzaldehyde, 4-hydroxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.244	110.72 ng/ul	2759840	Acenaphthene-d10	14.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	97
2			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	96
3			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	95
4			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	94
5			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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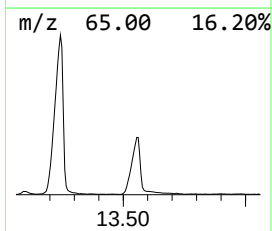
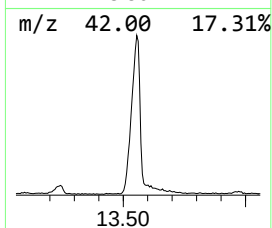
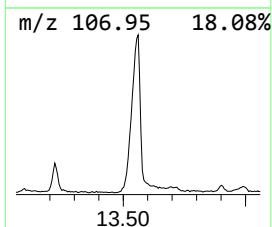
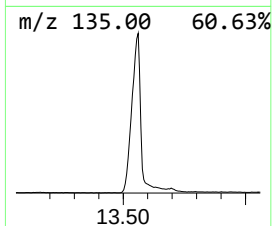
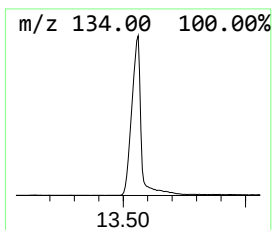
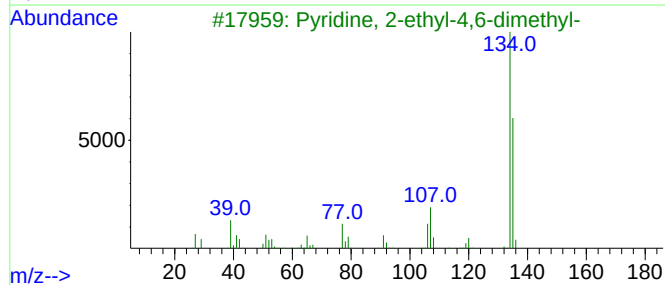
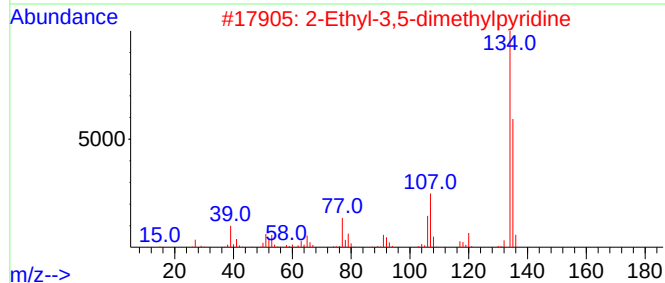
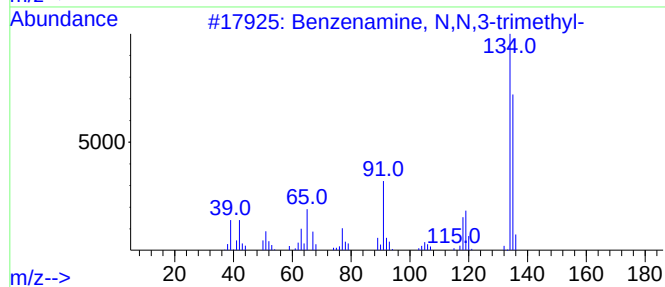
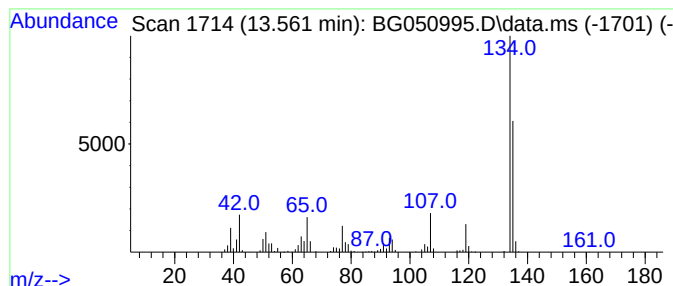
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 Benzenamine, N,N,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.561	109.84 ng/ul	2737930	Acenaphthene-d10	14.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, N,N,3-trimethyl-	135	C9H13N	000121-72-2	83
2			2-Ethyl-3,5-dimethylpyridine	135	C9H13N	001123-96-2	72
3			Pyridine, 2-ethyl-4,6-dimethyl-	135	C9H13N	001124-35-2	72
4			Pyridine, 2,6-diethyl-	135	C9H13N	000935-28-4	72
5			Benzenamine, N,N,4-trimethyl-	135	C9H13N	000099-97-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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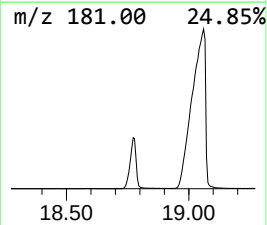
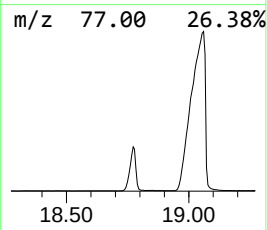
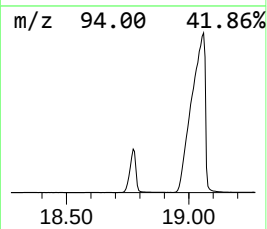
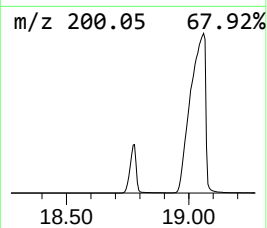
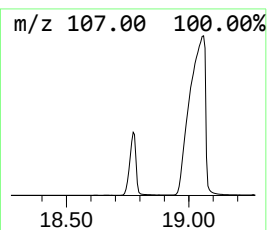
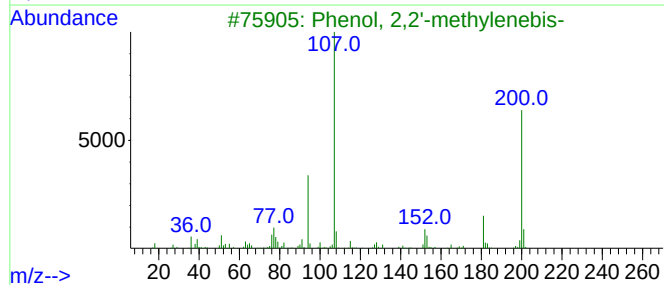
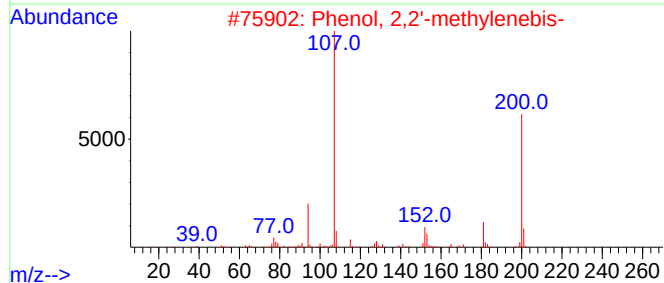
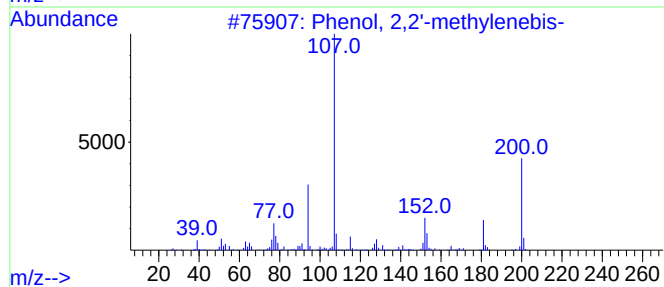
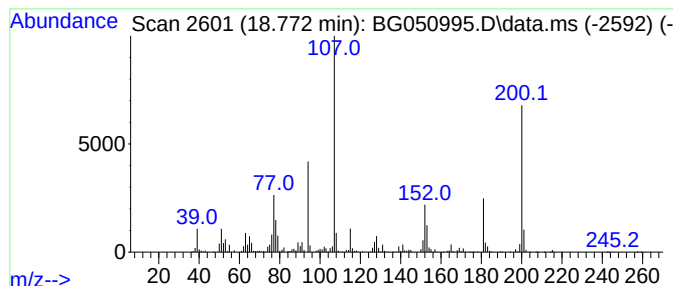
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 Phenol, 2,2'-methylenebis- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.772	243.55 ng/ul	6461020	Phenanthrene-d10	17.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	97
2			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	96
3			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	93
4			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	90
5			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

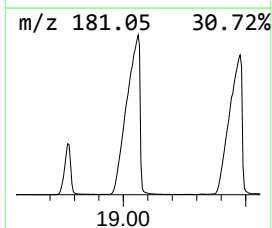
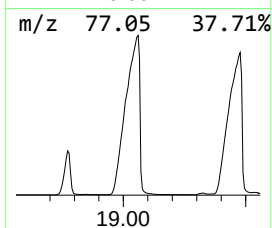
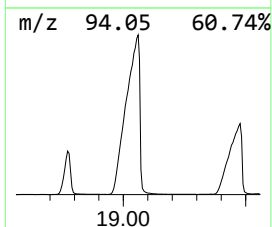
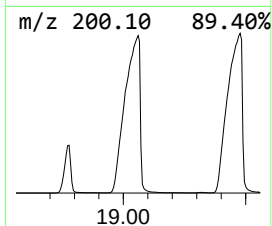
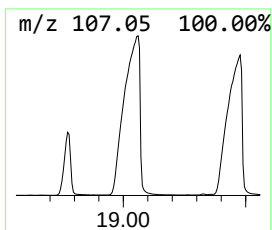
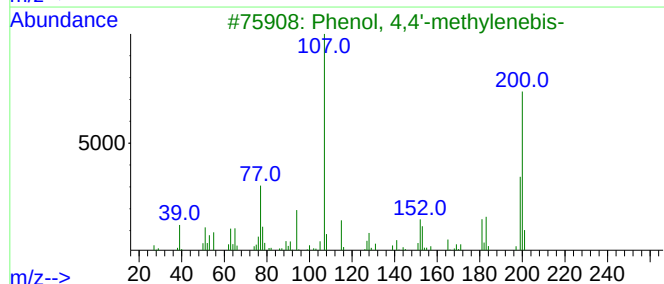
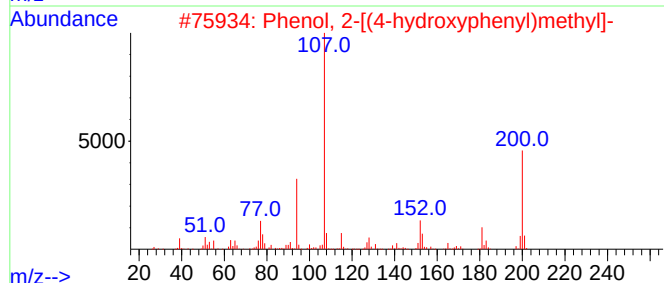
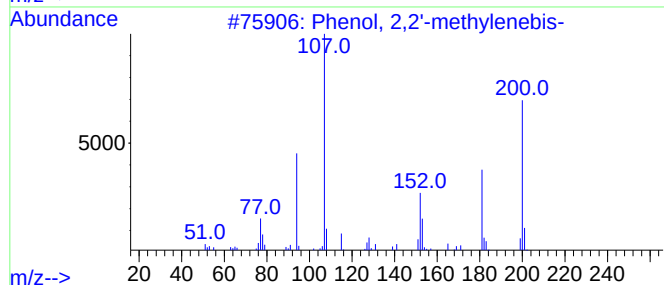
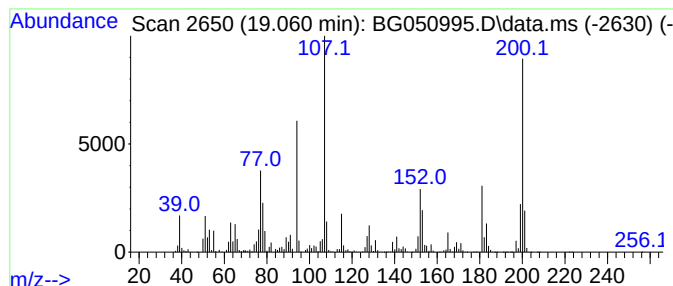
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.060	2072.51 ng/ul	54980900	Phenanthrene-d10	17.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	90
2			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	83
3			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	62
4			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	60
5			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

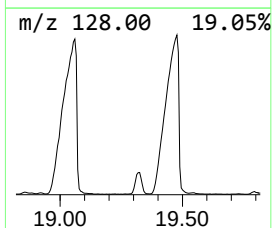
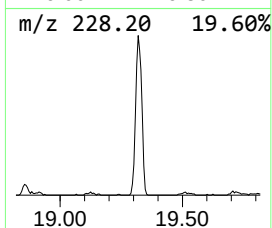
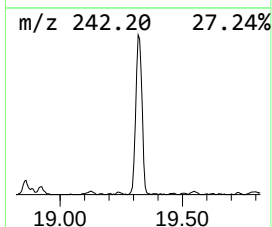
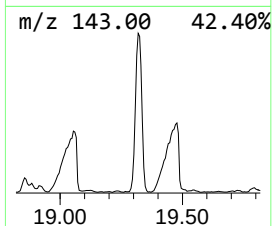
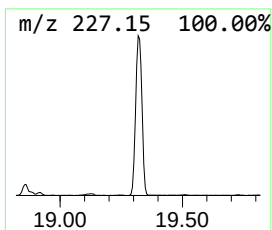
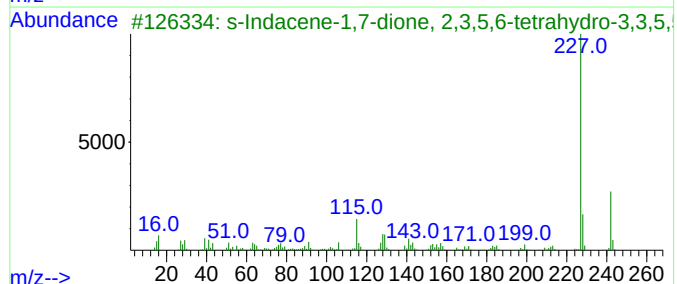
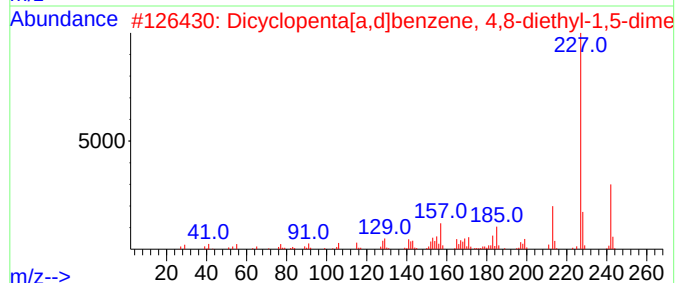
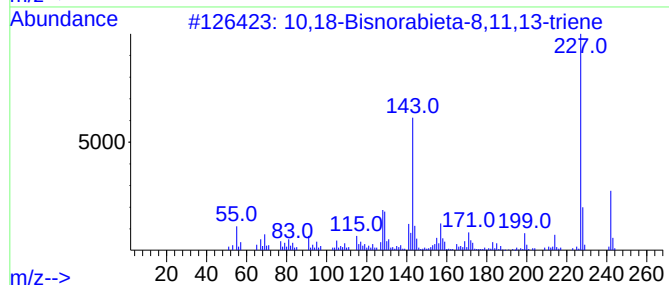
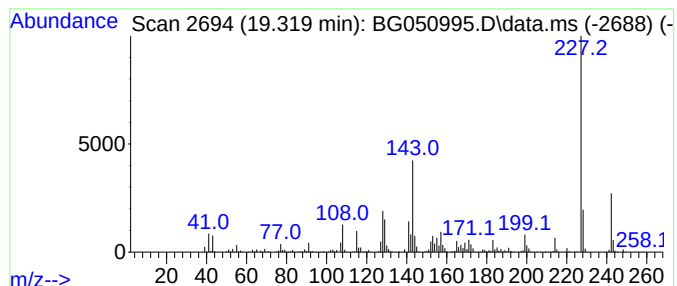
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 10,18-Bisnorabieta-8,11,13-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.319	51.78 ng/ul	1373650	Phenanthrene-d10	17.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	94		
2	Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	60		
3	s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	49		
4	Lapachol	242	C15H14O3	000084-79-7	49		
5	benzenesulfonamide, 4-[(diethyla...	242	C11H18N2O2S	1000396-54-0	46		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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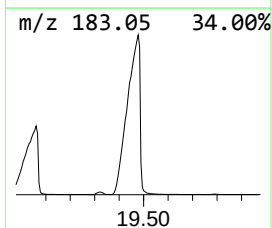
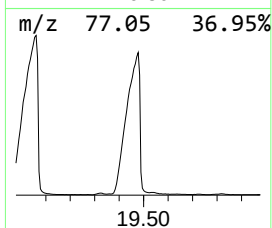
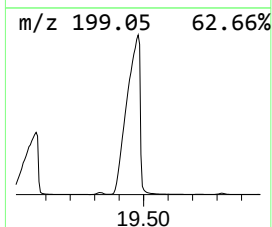
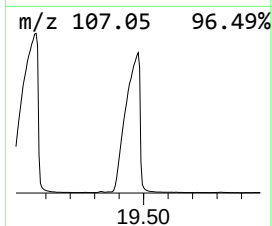
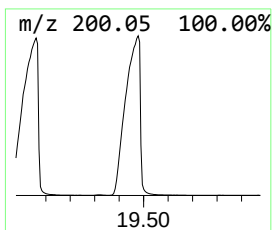
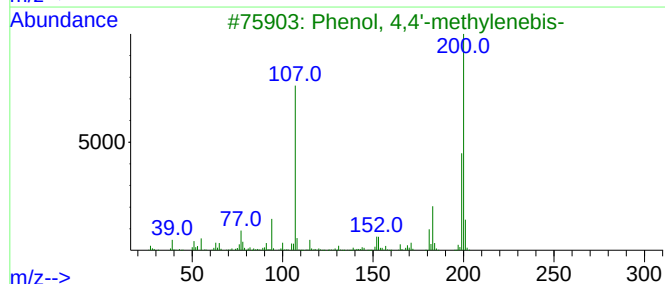
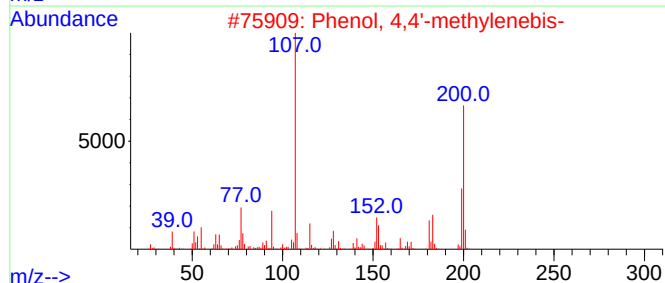
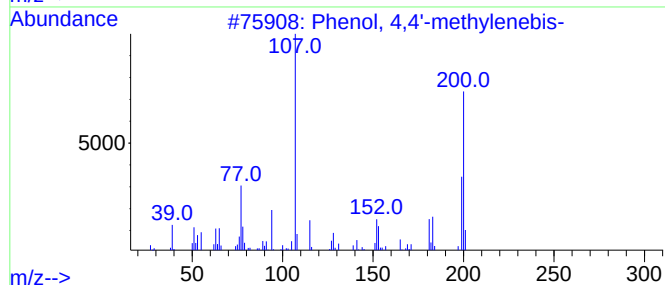
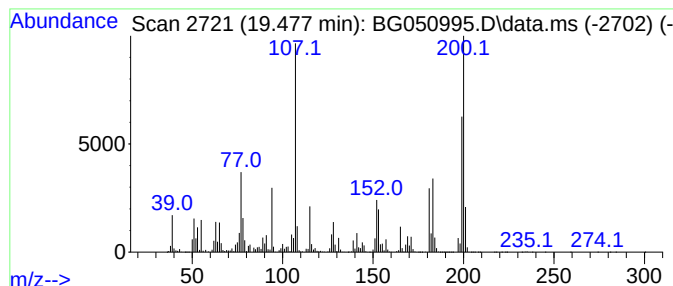
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 Phenol, 4,4'-methylenebis- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.477	1899.94 ng/ul	50402900	Phenanthrene-d10	17.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	95
2			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	93
3			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	81
4			Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	68
5			Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

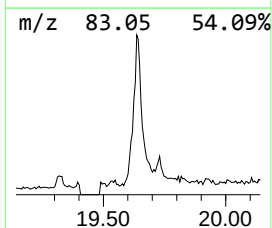
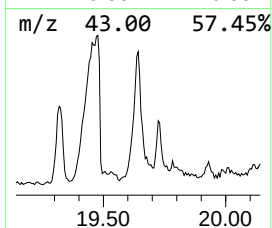
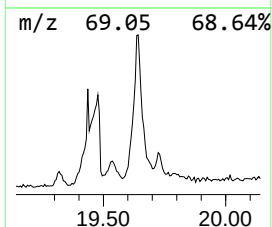
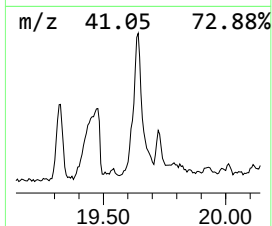
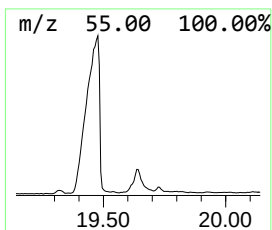
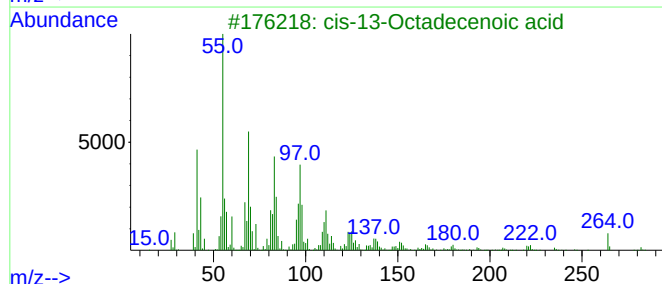
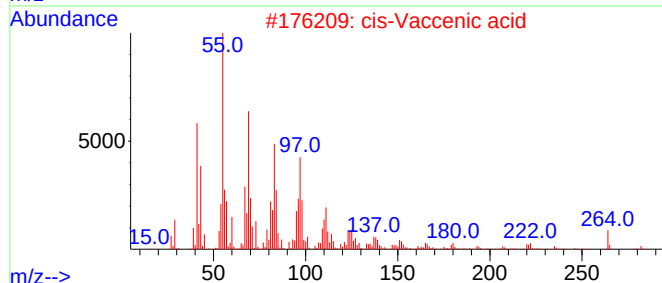
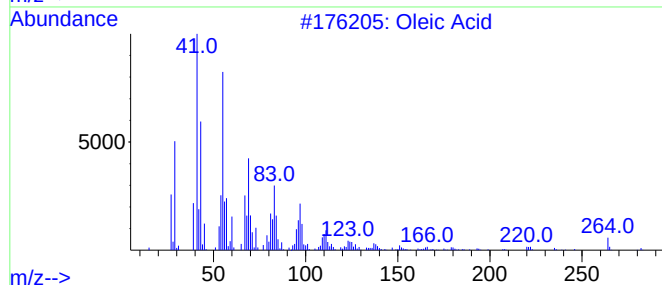
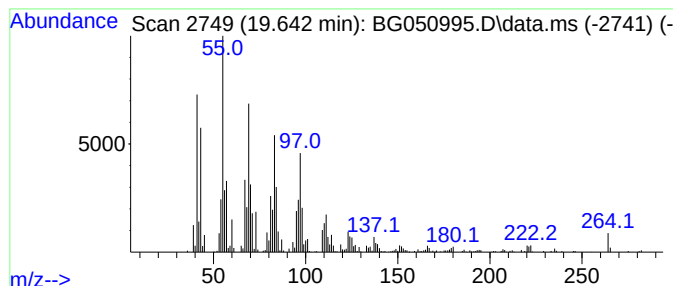
Instrument :
BNA_G
ClientSampleId :
COV04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 Oleic Acid Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.642	30.12 ng/ul	799005	Phenanthrene-d10		17.603	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Oleic Acid		282	C18H34O2	000112-80-1	99
2	cis-Vaccenic acid		282	C18H34O2	000506-17-2	94
3	cis-13-Octadecenoic acid		282	C18H34O2	013126-39-1	92
4	9-Octadecenoic acid, (E)-		282	C18H34O2	000112-79-8	91
5	Palmitoleic acid		254	C16H30O2	000373-49-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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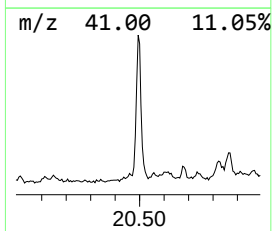
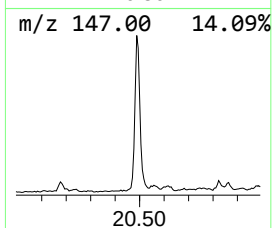
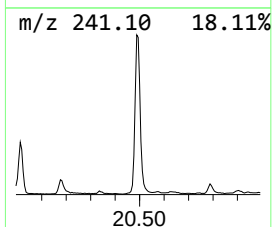
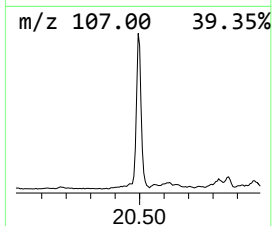
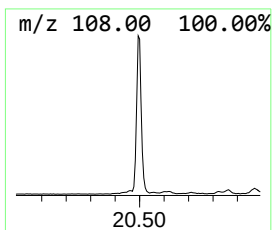
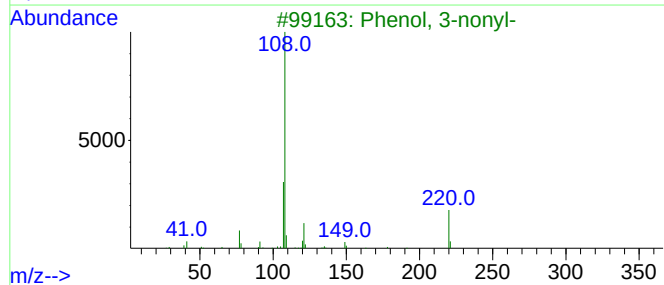
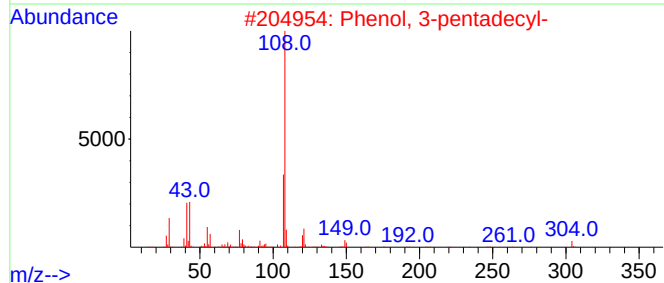
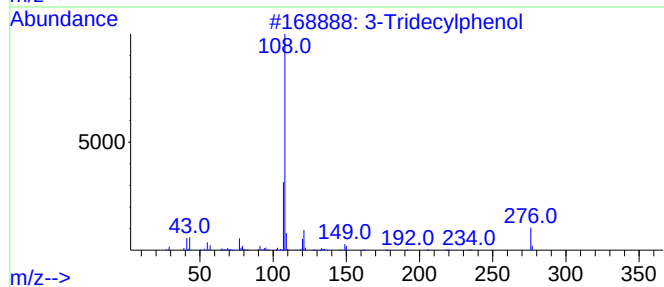
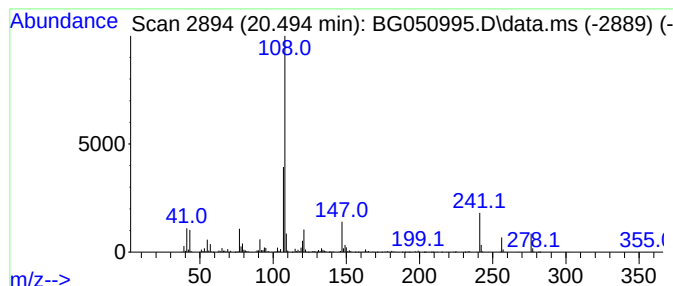
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 3-Tridecylphenol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.494	59.20 ng/ul	1657660	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Tridecylphenol	276	C19H32O	072424-02-3	93
2			Phenol, 3-pentadecyl-	304	C21H36O	000501-24-6	64
3			Phenol, 3-nonyl-	220	C15H24O	000139-84-4	64
4			Phenol, 3-undecyl-	248	C17H28O	020056-72-8	64
5			Acetic acid, 4-methylphenyl ester	150	C9H10O2	000140-39-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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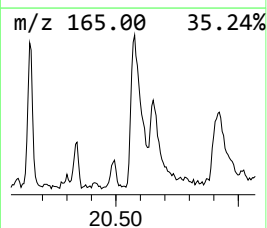
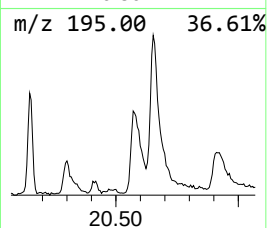
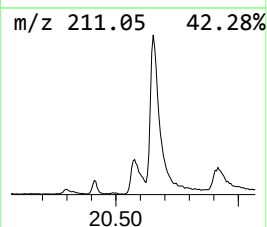
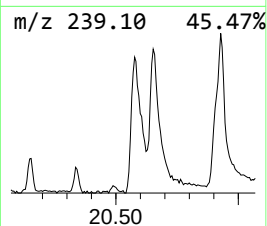
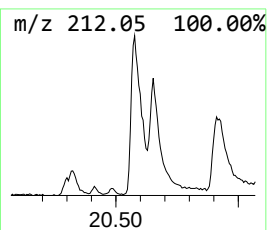
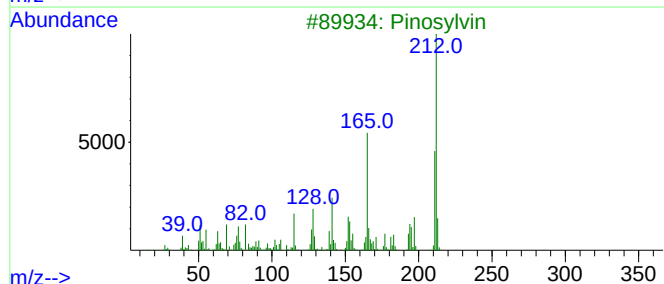
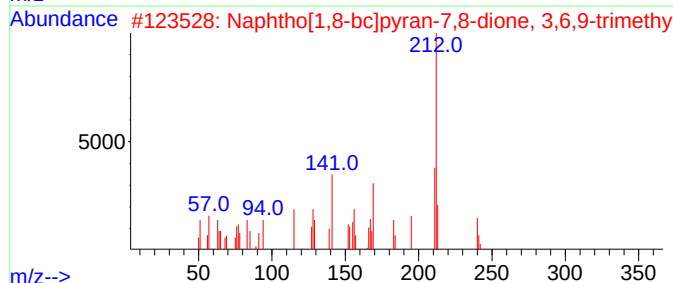
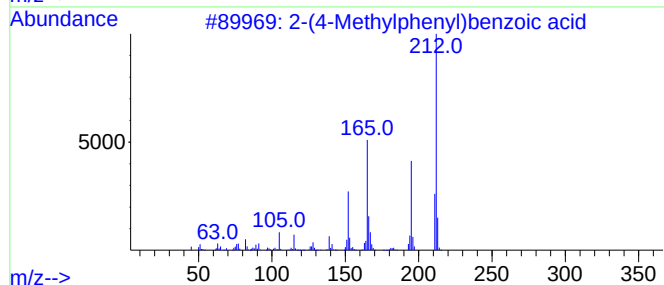
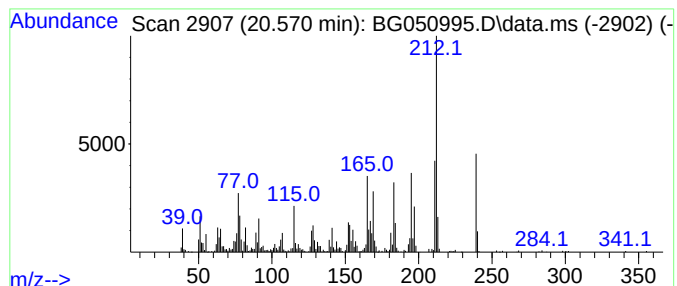
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 2-(4-Methylphenyl)benzoic acid Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.570	30.34 ng/ul	849519	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-(4-Methylphenyl)benzoic acid	212	C14H12O2	1000305-27-8	64
2			Naphtho[1,8-bc]pyran-7,8-dione, ...	240	C15H12O3	005090-88-0	58
3			Pinosylvins	212	C14H12O2	022139-77-1	55
4			2,4'-Dihydroxy-stilbene	212	C14H12O2	1000147-73-5	55
5			9H-Xanthen-9-one, 2-hydroxy-	212	C13H8O3	001915-98-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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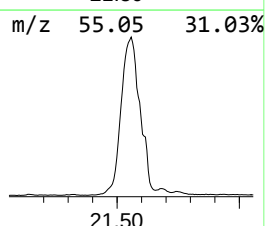
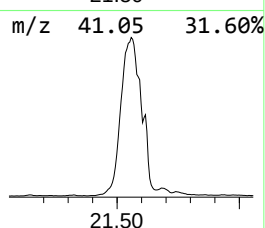
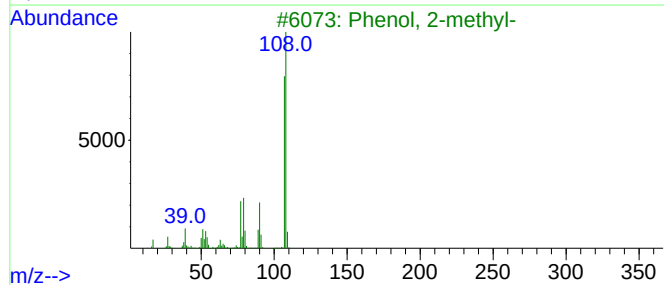
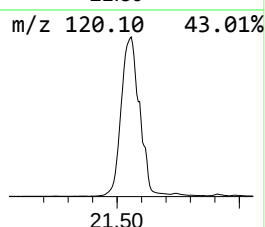
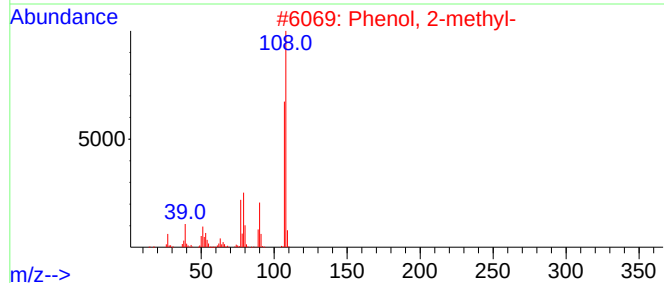
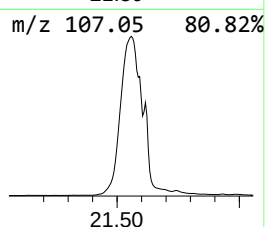
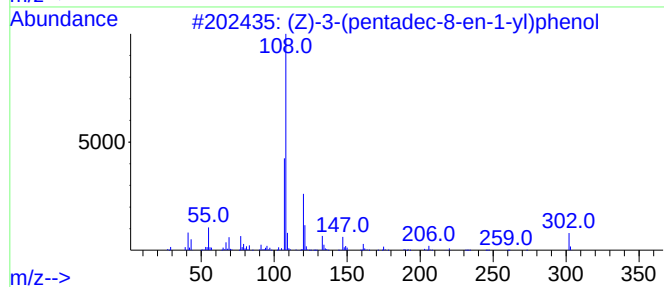
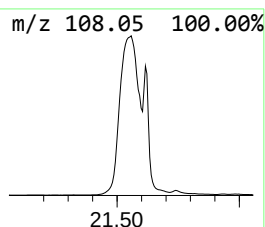
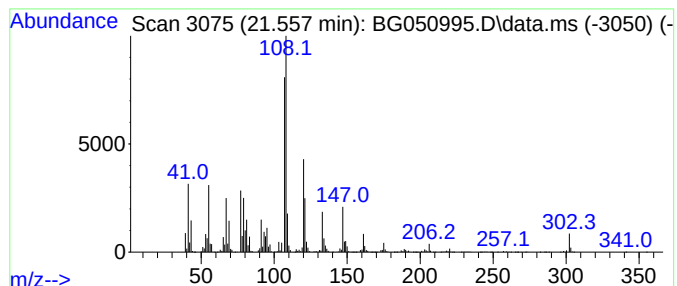
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 (Z)-3-(pentadec-8-en-1-yl)p... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.557	2095.44 ng/ul	58671700	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-3-(pentadec-8-en-1-yl)phenol	302	C21H34O	000501-26-8	53		
2	Phenol, 2-methyl-	108	C7H8O	000095-48-7	52		
3	Phenol, 2-methyl-	108	C7H8O	000095-48-7	52		
4	p-Cresol	108	C7H8O	000106-44-5	52		
5	Phenol, 3-methyl-	108	C7H8O	000108-39-4	52		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

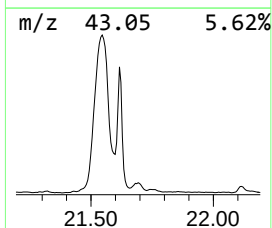
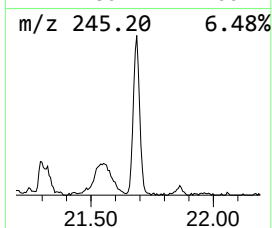
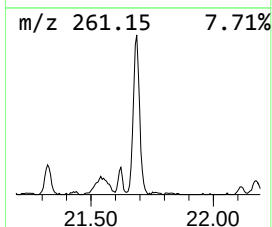
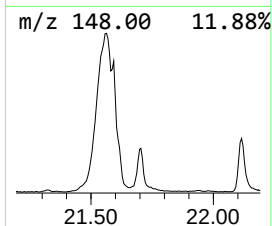
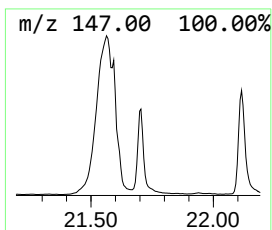
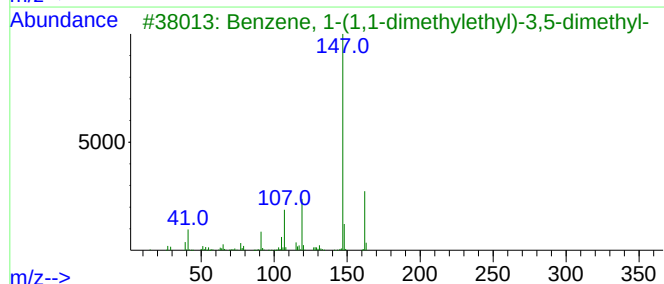
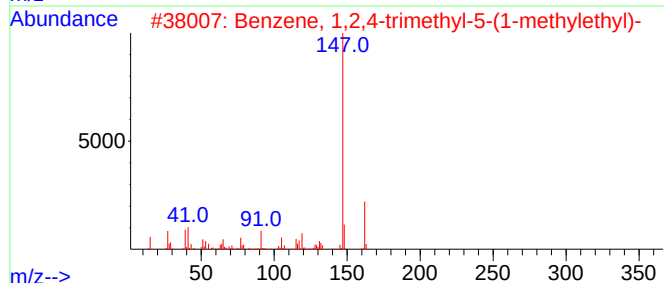
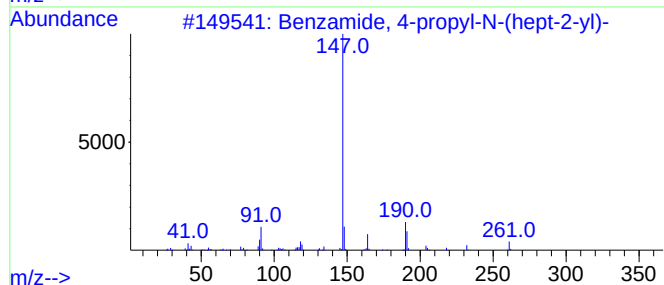
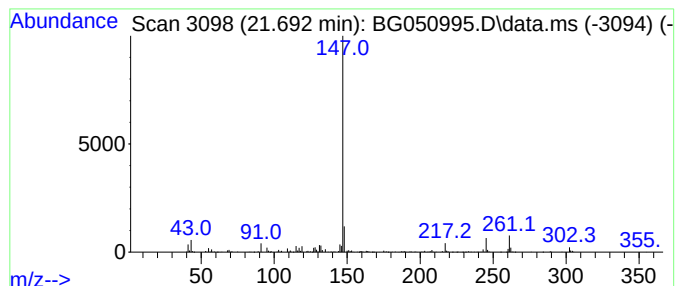
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 72 Benzamide, 4-propyl-N-(hept... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.692	38.54 ng/ul	1079060	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, 4-propyl-N-(hept-2-yl)-	261	C17H27NO	1000439-17-0	64
2			Benzene, 1,2,4-trimethyl-5-(1-me...	162	C12H18	010222-95-4	59
3			Benzene, 1-(1,1-dimethylethyl)-3...	162	C12H18	000098-19-1	59
4			Pentanoic acid, 5-oxo-5-[(5,6,7,...	261	C15H19NO3	1000337-45-4	50
5			2,2-Dibromo-1-(2,4,6-trimethylph...	318	C11H12Br2O	329349-15-7	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

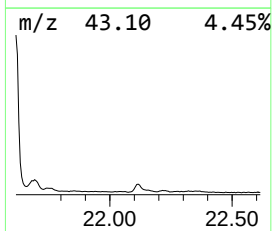
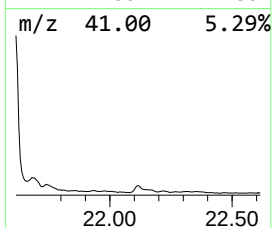
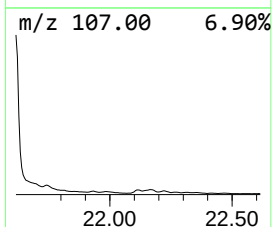
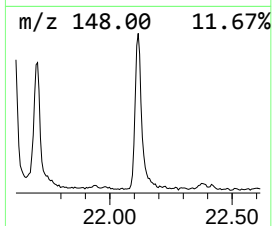
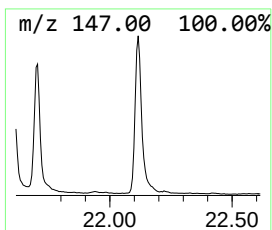
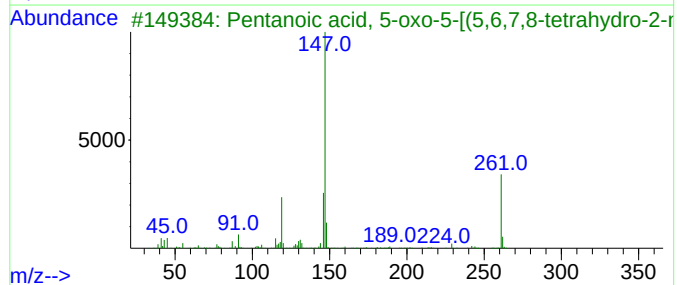
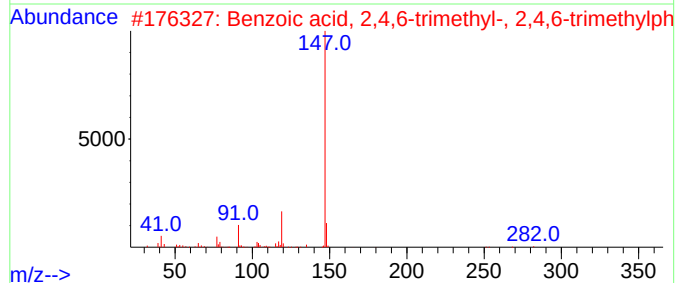
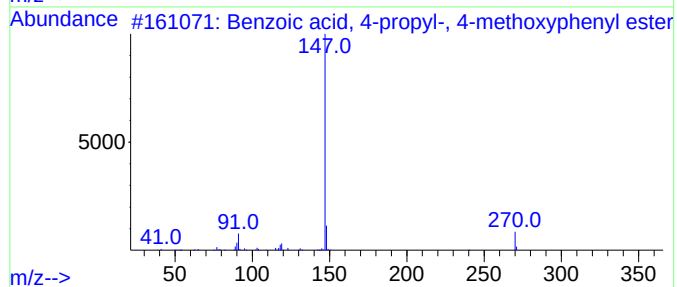
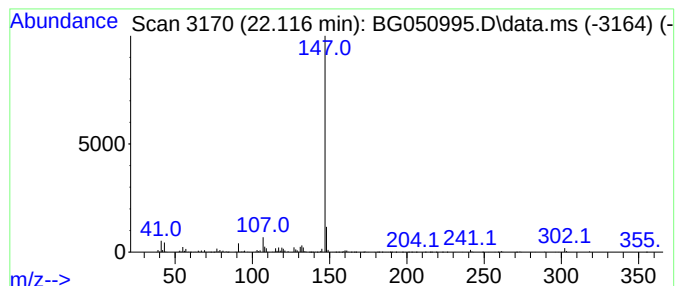
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 73 Benzoic acid, 4-propyl-, 4-... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.116	49.53 ng/ul	1386880	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 4-propyl-, 4-metho...	270	C17H18O3	050649-61-1	72
2			Benzoic acid, 2,4,6-trimethyl-, ...	282	C19H22O2	001504-38-7	64
3			Pentanoic acid, 5-oxo-5-[(5,6,7,...	261	C15H19NO3	1000337-45-4	64
4			1,3-dithiolan-2-one, 4-(iodometh...	274	C5H7IOS2	1010397-03-4	53
5			DL-Nornicotine, N-pentafluoropro...	294	C12H11F5N2O	1000448-18-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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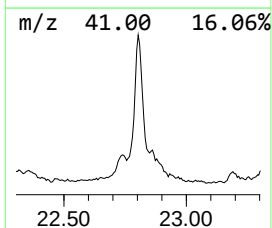
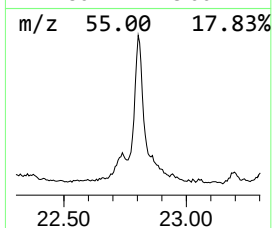
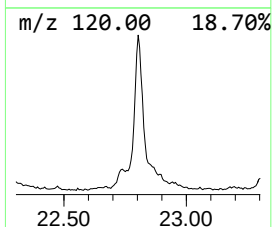
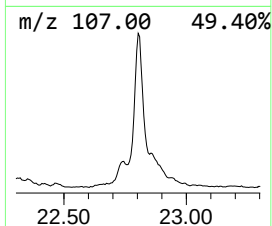
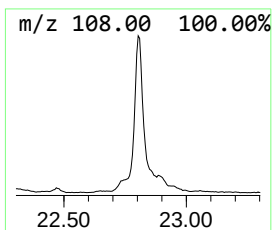
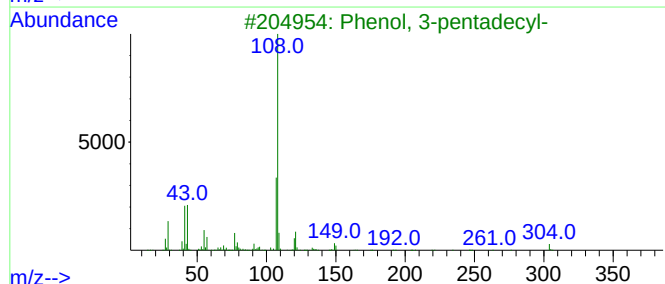
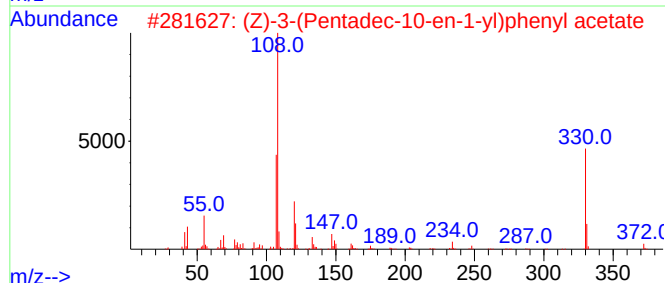
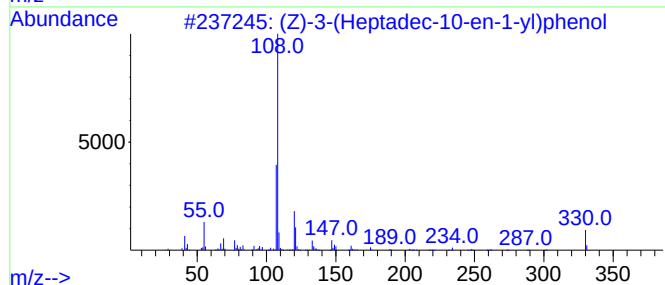
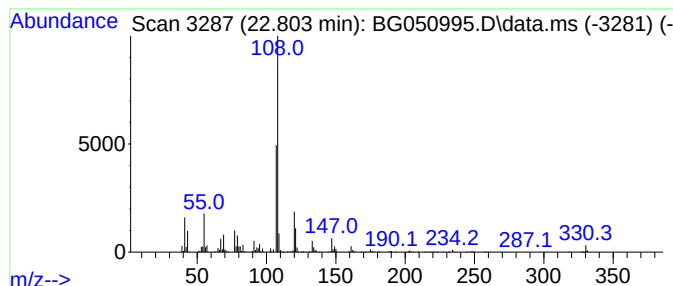
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 (Z)-3-(Heptadec-10-en-1-yl)... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.803	62.14 ng/ul	1740020	Chrysene-d12	21.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-3-(Heptadec-10-en-1-yl)phenol	330	C23H38O	111047-33-7	95		
2	(Z)-3-(Pentadec-10-en-1-yl)phenyl...	372	C25H40O2	454464-13-2	91		
3	Phenol, 3-pentadecyl-	304	C21H36O	000501-24-6	72		
4	(Z)-3-(pentadec-8-en-1-yl)phenol	302	C21H34O	000501-26-8	72		
5	Phenol, 2-methyl-	108	C7H8O	000095-48-7	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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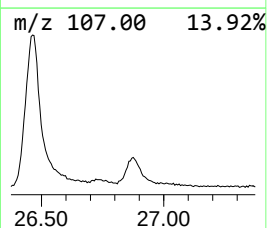
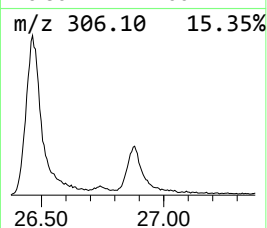
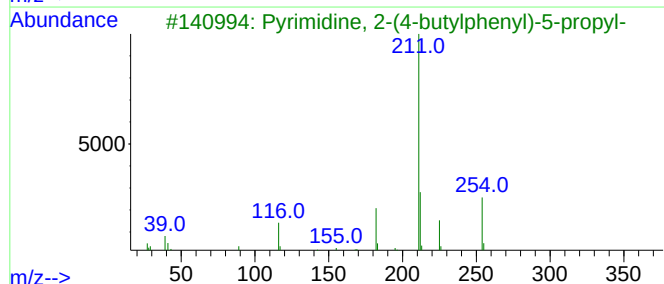
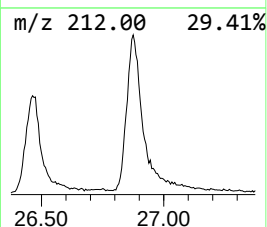
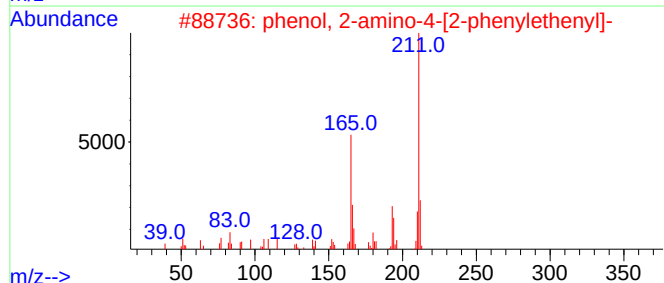
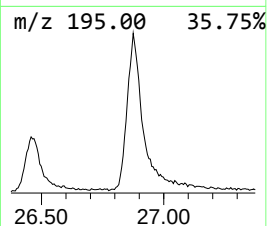
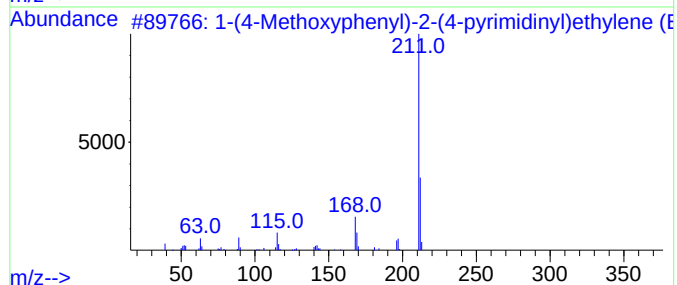
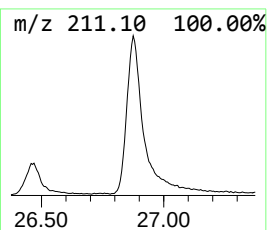
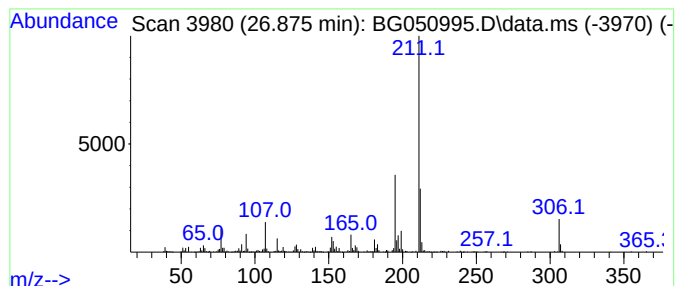
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 78 1-(4-Methoxyphenyl)-2-(4-py... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.875	46.03 ng/ul	1207590	Perylene-d12	25.312

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(4-Methoxyphenyl)-2-(4-pyrimidin...	212	C13H12N2O	163011-00-5	52
2			phenol, 2-amino-4-[2-phenylethen...	211	C14H13NO	1000396-15-8	47
3			Pyrimidine, 2-(4-butylphenyl)-5-...	254	C17H22N2	123740-94-3	38
4			Alnusone	294	C19H18O3	052330-11-7	37
5			1H-Benzo[d,E]phthalazine, 6-meth...	226	C14H14N2O	078378-22-0	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
Data File : BG050995.D
Acq On : 12 Nov 2021 10:58
Operator : CG/JU
Sample : M4615-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COV04

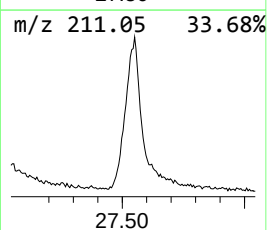
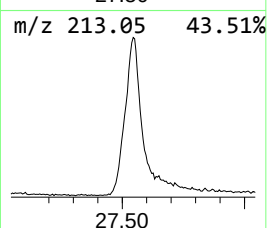
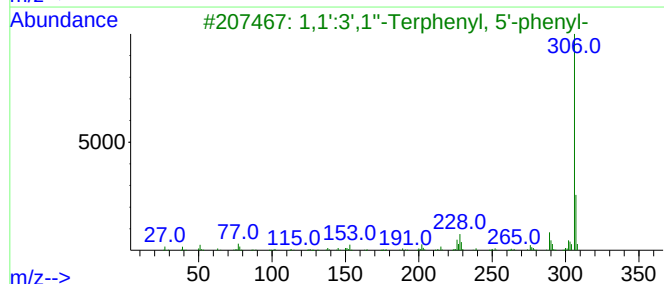
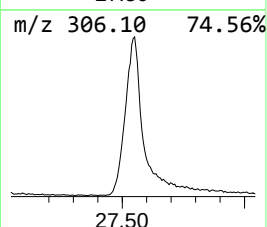
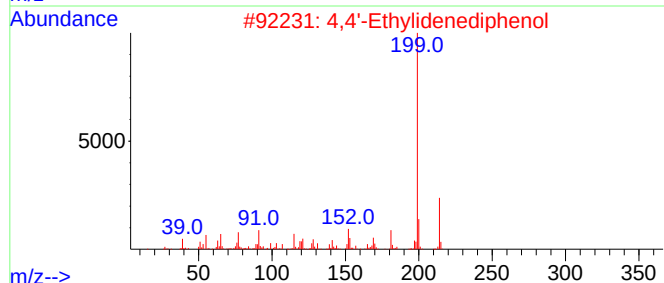
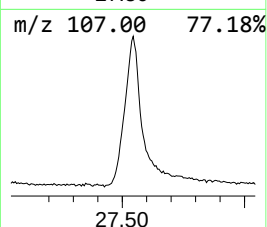
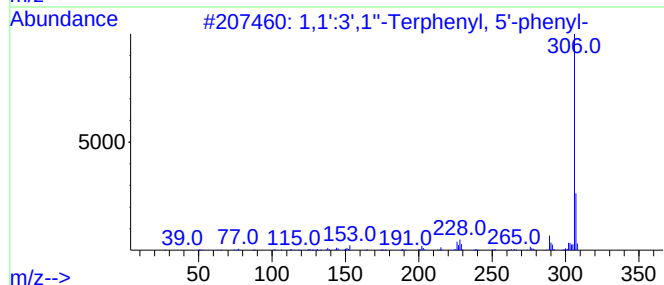
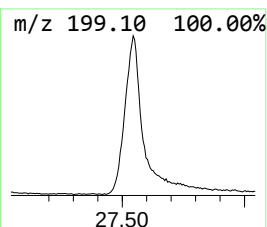
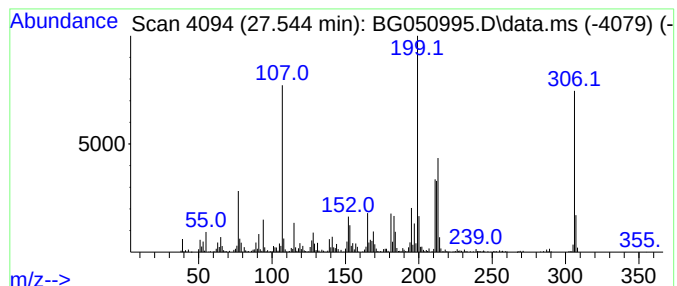
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 79 unknown-03 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.544	99.13 ng/ul	2600520	Perylene-d12	25.312

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1':3',1''-Terphenyl, 5'-phenyl-	306	C24H18	000612-71-5	25		
2	4,4'-Ethylidenediphenol	214	C14H14O2	002081-08-5	25		
3	1,1':3',1''-Terphenyl, 5'-phenyl-	306	C24H18	000612-71-5	25		
4	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	15		
5	1,1,2,2-Tetrakis(4-hydroxyphenyl...	398	C26H22O4	007727-33-5	15		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050995.D
 Acq On : 12 Nov 2021 10:58
 Operator : CG/JU
 Sample : M4615-05
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COV04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid, 1,...	3.526	223.3	ng/ul	12074300	1	8.261	1081190	20.0
(DEL) Alkane: C...	5.276	4.3	ng/ul	233397	1	8.261	1081190	20.0
(DEL) Alkane: C...	5.629	15.0	ng/ul	811622	1	8.261	1081190	20.0
(DEL) Alkane: S...	5.764	5.0	ng/ul	269593	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.005	7.7	ng/ul	413819	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.070	14.9	ng/ul	802580	1	8.261	1081190	20.0
(DEL) Alkane: S...	6.158	12.4	ng/ul	667742	1	8.261	1081190	20.0
(DEL) Alkane: S...	6.211	44.8	ng/ul	2421790	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.481	35.0	ng/ul	1894360	1	8.261	1081190	20.0
(DEL) Alkane: S...	6.587	7.6	ng/ul	412927	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.710	14.7	ng/ul	793039	1	8.261	1081190	20.0
(DEL) Alkane: S...	6.757	23.7	ng/ul	1280210	1	8.261	1081190	20.0
(DEL) Alkane: S...	6.810	4.7	ng/ul	252701	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.851	47.7	ng/ul	2580170	1	8.261	1081190	20.0
(DEL) Alkane: C...	6.963	11.6	ng/ul	625281	1	8.261	1081190	20.0
(DEL) Alkane: C...	7.033	7.2	ng/ul	386406	1	8.261	1081190	20.0
(DEL) Alkane: S...	7.098	15.2	ng/ul	819640	1	8.261	1081190	20.0
(DEL) Alkane: S...	7.157	8.6	ng/ul	464183	1	8.261	1081190	20.0
(DEL) Alkane: S...	7.198	35.3	ng/ul	1906650	1	8.261	1081190	20.0
(DEL) Alkane: S...	7.251	30.4	ng/ul	1645190	1	8.261	1081190	20.0
unknown-01	7.362	61.1	ng/ul	3303380	1	8.261	1081190	20.0
1,3,5-Triazine,...	7.750	51.1	ng/ul	2760500	1	8.261	1081190	20.0
Benzene, 1,2,3-...	7.915	33.3	ng/ul	1797300	1	8.261	1081190	20.0
(DEL) Alkane: S...	8.197	64.5	ng/ul	3485300	1	8.261	1081190	20.0
(DEL) Alkane: S...	8.473	16.7	ng/ul	904580	1	8.261	1081190	20.0
(DEL) Alkane: C...	8.520	49.8	ng/ul	2693930	1	8.261	1081190	20.0
(DEL) Alkane: C...	8.573	12.1	ng/ul	653201	1	8.261	1081190	20.0
(DEL) Alkane: S...	8.643	5.3	ng/ul	288837	1	8.261	1081190	20.0
unknown-02	8.790	55.5	ng/ul	3002750	1	8.261	1081190	20.0
(DEL) Alkane: S...	8.878	39.2	ng/ul	2120690	1	8.261	1081190	20.0
(DEL) Alkane: S...	9.301	7.8	ng/ul	422680	1	8.261	1081190	20.0
Benzene, 4-ethy...	9.342	29.8	ng/ul	1608920	1	8.261	1081190	20.0
(DEL) Alkane: C...	9.554	8.0	ng/ul	433099	1	8.261	1081190	20.0
(DEL) Alkane: S...	9.695	40.0	ng/ul	661174	2	11.064	330429	20.0
(DEL) Alkane: S...	9.865	36.0	ng/ul	594514	2	11.064	330429	20.0
o-Cymene	10.447	35.2	ng/ul	581320	2	11.064	330429	20.0
(DEL) Alkane: S...	10.993	49.8	ng/ul	823322	2	11.064	330429	20.0
Benzenamine, N,...	11.634	231.6	ng/ul	3826050	2	11.064	330429	20.0
Methenamine	11.851	1001.7	ng/ul	16549600	2	11.064	330429	20.0
4(1H)-Pyrimidin...	12.074	324.3	ng/ul	5358150	2	11.064	330429	20.0
Phenol, p-tert-...	12.398	76.7	ng/ul	1266790	2	11.064	330429	20.0
Benzaldehyde, 4...	13.244	110.7	ng/ul	2759840	3	14.859	498513	20.0
Benzenamine, N,...	13.561	109.8	ng/ul	2737930	3	14.859	498513	20.0
Phenol, 2,2'-me...	18.772	243.6	ng/ul	6461020	4	17.603	530573	20.0
Phenol, 2-[(4-h...	19.060	2072.5	ng/ul	54980900	4	17.603	530573	20.0
10,18-Bisnorabi...	19.319	51.8	ng/ul	1373650	4	17.603	530573	20.0
Phenol, 4,4'-me...	19.477	1899.9	ng/ul	50402900	4	17.603	530573	20.0
Oleic Acid	19.642	30.1	ng/ul	799005	4	17.603	530573	20.0
3-Tridecylphenol	20.494	59.2	ng/ul	1657660	5	21.910	559994	20.0
2-(4-Methylphen...	20.570	30.3	ng/ul	849519	5	21.910	559994	20.0
(Z)-3-(pentadec...	21.557	2095.4	ng/ul	58671700	5	21.910	559994	20.0
Benzamide, 4-pr...	21.692	38.5	ng/ul	1079060	5	21.910	559994	20.0

Benzoic acid, 4...	22.116	49.5	ng/ul	1386880	5	21.910	559994	20.0
(Z)-3-(Heptadec...	22.803	62.1	ng/ul	1740020	5	21.910	559994	20.0
1-(4-Methoxyphe...	26.875	46.0	ng/ul	1207590	6	25.312	524682	20.0
unknown-03	27.544	99.1	ng/ul	2600520	6	25.312	524682	20.0

Instrument :

BNA_G

ClientSampleId :

C0V04