

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051789.D
 Acq On : 23 Dec 2021 17:44
 Operator : CG/JU
 Sample : M5215-04
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLD6

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-BG121421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051789.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.747	203	214	224	rVB	313150	704441	0.57%	0.074%
2	5.164	276	285	295	rBV	251836	626512	0.51%	0.066%
3	5.440	327	332	348	rVB2	403430	1137962	0.92%	0.120%
4	5.616	352	362	368	rBV4	212717	859880	0.70%	0.091%
5	5.781	368	390	401	rBV	11122915	44576862	36.23%	4.696%
6	5.975	401	423	429	rVB3	23706904	123050761	100.00%	12.962%
7	6.086	437	442	446	rVB	483389	725524	0.59%	0.076%
8	6.151	446	453	461	rBV4	730064	2144063	1.74%	0.226%
9	6.233	461	467	469	rBV2	919283	1822462	1.48%	0.192%
10	6.304	469	479	487	rVB2	25741976	61984942	50.37%	6.530%
11	6.380	487	492	499	rBV3	495253	921625	0.75%	0.097%
12	6.451	499	504	513	rVB3	340611	774969	0.63%	0.082%
13	6.550	513	521	534	rBV6	2917497	9715193	7.90%	1.023%
14	6.668	534	541	550	rVB3	1763479	4690988	3.81%	0.494%
15	6.768	550	558	565	rBV6	3562229	10190953	8.28%	1.074%
16	6.844	565	571	577	rBV	2998809	6195784	5.04%	0.653%
17	6.926	577	585	600	rVB4	4993410	18220675	14.81%	1.919%
18	7.044	601	605	611	rVB2	621432	1187757	0.97%	0.125%
19	7.108	611	616	621	rBV3	944508	1947653	1.58%	0.205%
20	7.191	622	630	636	rBV4	1646486	5162164	4.20%	0.544%
21	7.273	637	644	650	rBV	5874763	16730662	13.60%	1.762%
22	7.379	656	662	665	rBV	9775825	19772266	16.07%	2.083%
23	7.473	672	678	684	rVB3	11518313	26172047	21.27%	2.757%
24	7.549	684	691	697	rVB	10821086	24866907	20.21%	2.620%
25	7.649	697	708	711	rBV5	1901619	5183769	4.21%	0.546%
26	7.708	711	718	723	rBV	7784339	14635591	11.89%	1.542%
27	7.813	729	736	743	rBV3	1539591	4845361	3.94%	0.510%
28	7.984	754	765	767	rBV3	25199636	65985093	53.62%	6.951%
29	8.072	778	780	784	rVB2	1375281	1684680	1.37%	0.177%
30	8.125	785	789	792	rBV4	1165107	2002877	1.63%	0.211%
31	8.166	792	796	799	rBV4	835535	1432325	1.16%	0.151%
32	8.236	804	808	811	rBV	1528765	2786067	2.26%	0.293%
33	8.336	816	825	829	rVV	10039337	22183241	18.03%	2.337%
34	8.413	829	838	840	rVV3	2615026	6884663	5.59%	0.725%
35	8.471	840	848	852	rVB	13062575	25546724	20.76%	2.691%
36	8.530	852	858	862	rBV3	3749609	6359029	5.17%	0.670%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Title : SVOA CALIBRATION

37	8.589	862	868	872	rVV	5238104	10546442	8.57%	1.111%
38	8.642	872	877	881	rVV3	5816722	9779200	7.95%	1.030%
39	8.718	881	890	895	rVB3	4674092	11791504	9.58%	1.242%
40	8.771	895	899	903	rBV4	1190617	2193920	1.78%	0.231%
41	8.818	903	907	910	rBV	1573120	2633762	2.14%	0.277%
42	8.883	910	918	923	rVV4	11903611	27654087	22.47%	2.913%
43	8.935	924	927	930	rVV	4144599	5530304	4.49%	0.583%
44	8.988	930	936	949	rVB3	13667934	38314825	31.14%	4.036%
45	9.112	949	957	960	rBV	5809076	11639315	9.46%	1.226%
46	9.147	960	963	966	rVB	2454293	3178887	2.58%	0.335%
47	9.200	966	972	975	rBV2	2905973	5315659	4.32%	0.560%
48	9.288	983	987	991	rVB	2142143	3385379	2.75%	0.357%
49	9.347	991	997	1004	rVB	4232374	8759405	7.12%	0.923%
50	9.452	1004	1015	1023	rBV3	9817672	22561019	18.33%	2.377%
51	9.540	1023	1030	1031	rVV3	5171938	8171535	6.64%	0.861%
52	9.617	1033	1043	1047	rVB2	17891185	51968417	42.23%	5.474%
53	9.699	1048	1057	1066	rBV8	4194276	13937625	11.33%	1.468%
54	9.793	1066	1073	1075	rVV2	2573514	5409227	4.40%	0.570%
55	9.822	1076	1078	1086	rVV6	3406814	5960181	4.84%	0.628%
56	9.916	1087	1094	1099	rVV4	2170495	4612840	3.75%	0.486%
57	9.975	1099	1104	1109	rVV	4800319	8981794	7.30%	0.946%
58	10.040	1109	1115	1119	rVV2	4671321	8884141	7.22%	0.936%
59	10.093	1119	1124	1131	rVB3	2652863	5100141	4.14%	0.537%
60	10.210	1137	1144	1148	rBV4	1745000	4205984	3.42%	0.443%
61	10.269	1148	1154	1158	rVV3	3755442	7269594	5.91%	0.766%
62	10.334	1159	1165	1170	rVV2	3874870	7856048	6.38%	0.828%
63	10.404	1170	1177	1183	rVV2	2445764	6793807	5.52%	0.716%
64	10.474	1183	1189	1195	rVV3	2648710	5437024	4.42%	0.573%
65	10.551	1195	1202	1207	rVV3	5671389	11838083	9.62%	1.247%
66	10.604	1207	1211	1215	rVV3	1653764	2947366	2.40%	0.310%
67	10.657	1215	1220	1223	rVV	1869741	3110352	2.53%	0.328%
68	10.715	1223	1230	1235	rVV4	1246497	3462974	2.81%	0.365%
69	10.774	1236	1240	1245	rVV	1147896	2064690	1.68%	0.217%
70	10.833	1245	1250	1259	rVB2	852508	2058887	1.67%	0.217%
71	10.915	1259	1264	1268	rBV6	286982	520970	0.42%	0.055%
72	11.027	1279	1283	1289	rVV4	806026	1736447	1.41%	0.183%
73	11.115	1289	1298	1303	rVV	8214908	16606713	13.50%	1.749%
74	11.197	1303	1312	1317	rVV2	3140596	6883918	5.59%	0.725%
75	11.285	1317	1327	1333	rVB2	2612704	5401326	4.39%	0.569%
76	11.350	1333	1338	1344	rBV2	809739	1485221	1.21%	0.156%

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77	11.503	1358	1364	1369	rBV6	423428	888808	0.72%	0.094%
78	11.831	1414	1420	1425	rBV3	958850	1857152	1.51%	0.196%
79	11.949	1435	1440	1447	rVB4	781405	1388923	1.13%	0.146%
80	12.025	1447	1453	1461	rVV	1130654	2074378	1.69%	0.219%
81	12.131	1465	1471	1476	rVV	2082508	3819002	3.10%	0.402%
82	12.172	1476	1478	1482	rVV2	442395	636452	0.52%	0.067%
83	12.225	1482	1487	1496	rVB5	451472	905133	0.74%	0.095%
84	12.525	1531	1538	1547	rVB	5083777	8492812	6.90%	0.895%
85	12.730	1567	1573	1575	rBV3	443855	792192	0.64%	0.083%
86	12.771	1575	1580	1587	rVB	2107621	3559590	2.89%	0.375%
87	12.983	1600	1616	1622	rBV	1143771	2383265	1.94%	0.251%
88	13.130	1637	1641	1645	rBV	403383	690663	0.56%	0.073%
89	13.171	1645	1648	1651	rBV	444532	619273	0.50%	0.065%
90	13.306	1667	1671	1680	rVB	562458	833779	0.68%	0.088%
91	13.394	1681	1686	1690	rVV	359366	530638	0.43%	0.056%
92	13.447	1690	1695	1702	rVB	930329	1372995	1.12%	0.145%
93	13.735	1737	1744	1752	rBV2	2518253	4146713	3.37%	0.437%
94	14.234	1824	1829	1834	rVV2	339608	613330	0.50%	0.065%
95	14.293	1834	1839	1844	rVB2	472937	790749	0.64%	0.083%
96	14.910	1934	1944	1951	rBV5	338075	1036566	0.84%	0.109%
97	15.709	2076	2080	2083	rBV	394167	527911	0.43%	0.056%
98	16.631	2234	2237	2242	rVB	484420	623194	0.51%	0.066%
99	25.231	3693	3701	3711	rVB2	168099	509287	0.41%	0.054%
100	31.499	4757	4768	4788	rVB2	90971	502085	0.41%	0.053%

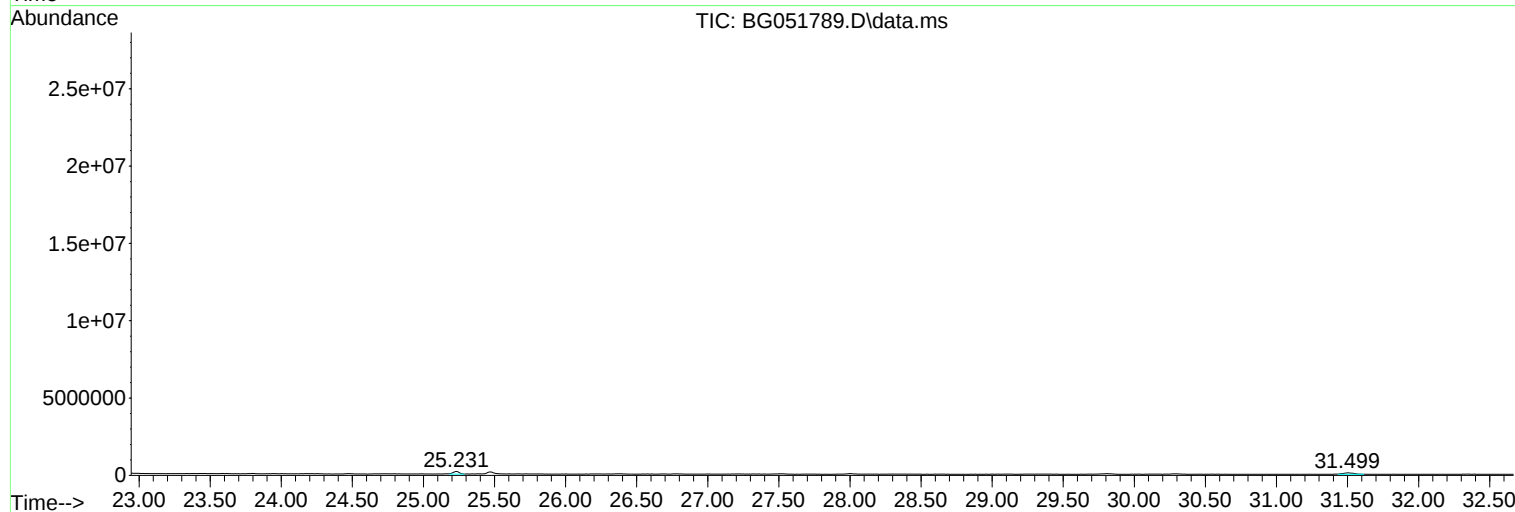
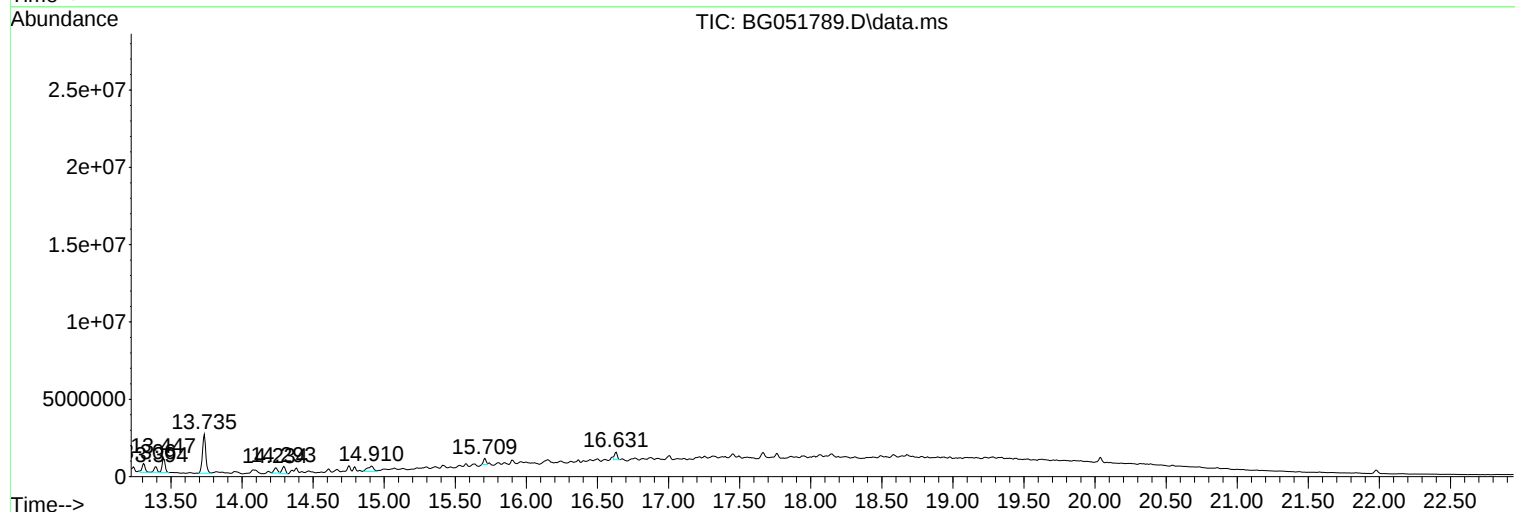
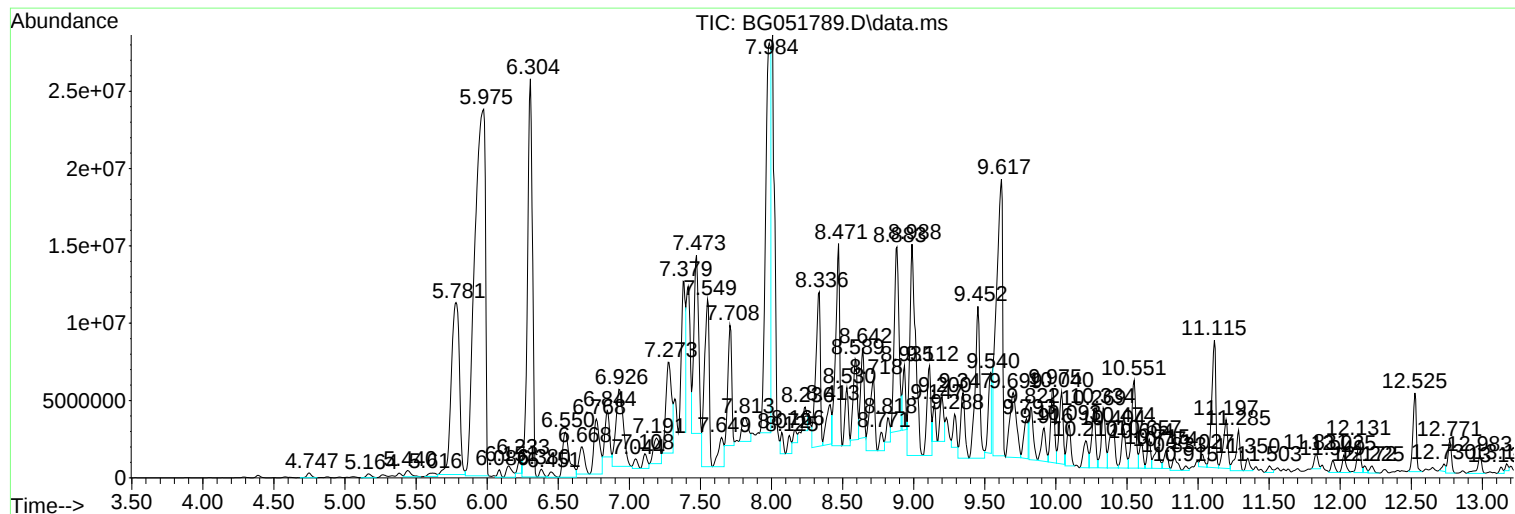
Sum of corrected areas: 949296375

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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



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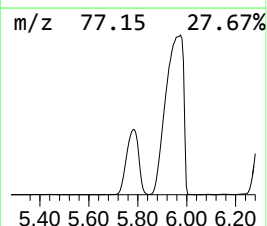
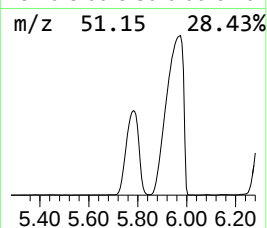
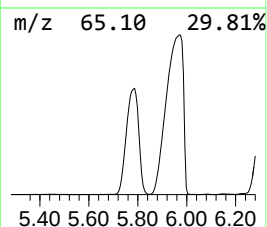
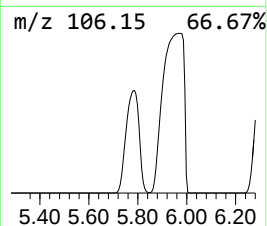
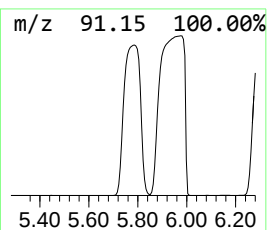
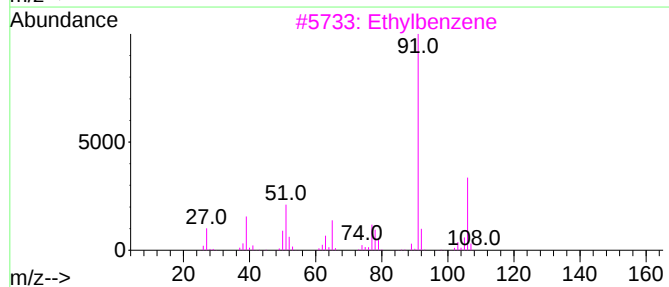
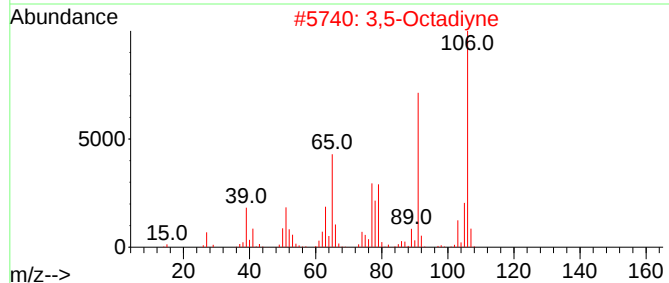
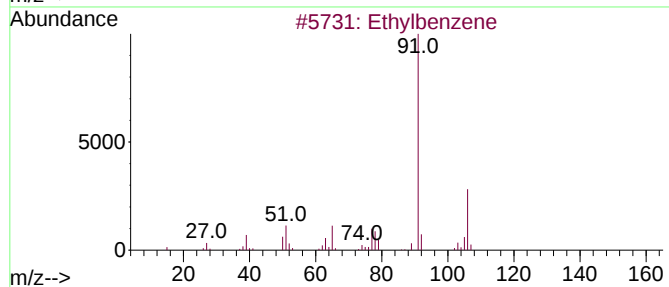
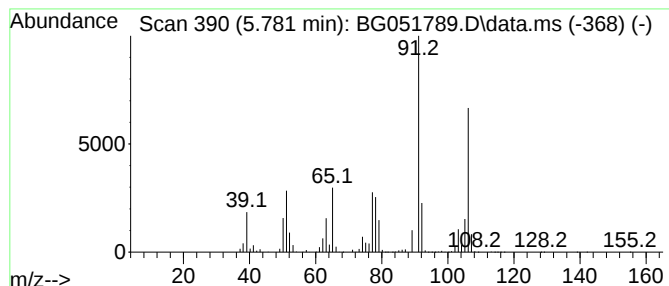
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethylbenzene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.781	40.19 ng/ul	44576900	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylbenzene	106	C8H10	000100-41-4	93
2			3,5-Octadiyne	106	C8H10	016387-70-5	76
3			Ethylbenzene	106	C8H10	000100-41-4	76
4			Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	58
5			Ethylbenzene	106	C8H10	000100-41-4	58



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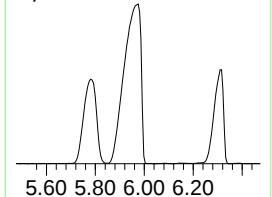
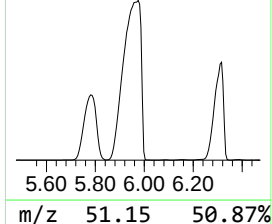
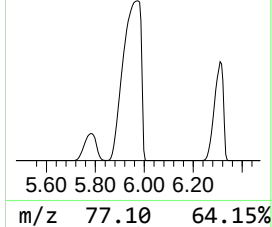
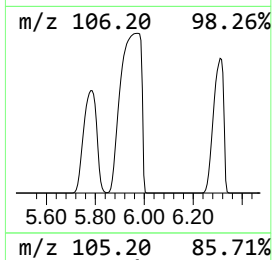
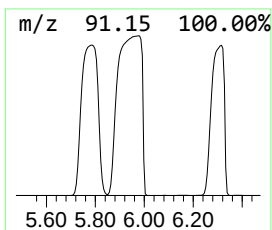
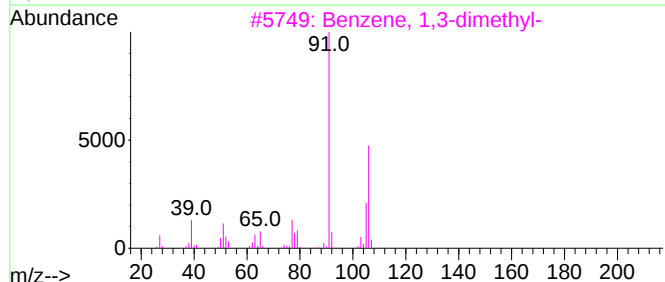
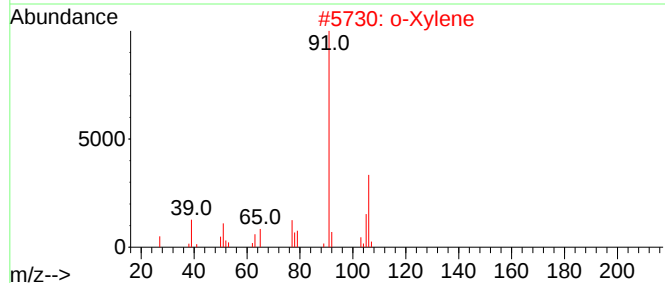
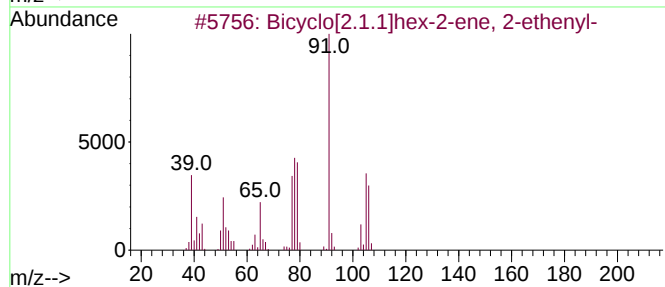
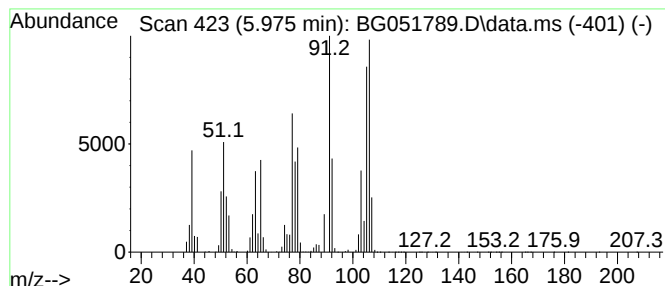
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Bicyclo[2.1.1]hex-2-ene, 2-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.975	110.94 ng/ul	123051000	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.1.1]hex-2-ene, 2-ethenyl-	106	C8H10	128600-91-9	64
2			o-Xylene	106	C8H10	000095-47-6	53
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	50
4			o-Xylene	106	C8H10	000095-47-6	50
5			5-Methylidenetricyclo[4.1.0.0(2,...	106	C8H10	1000436-95-4	49



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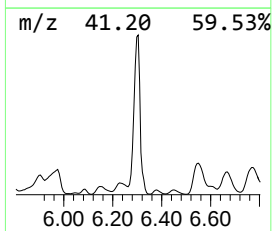
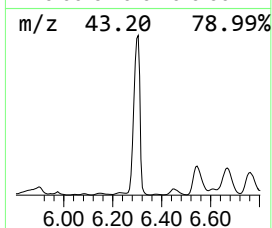
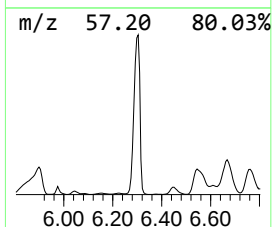
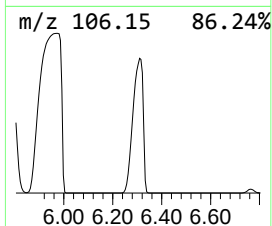
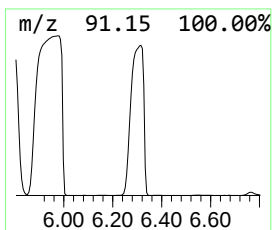
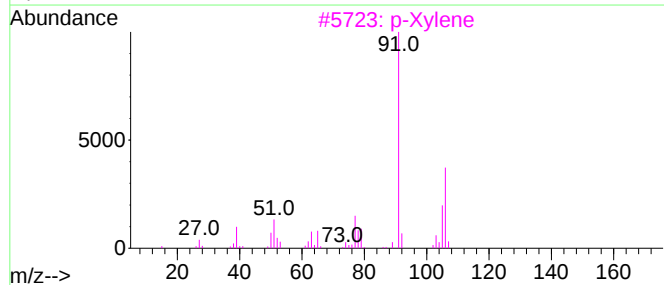
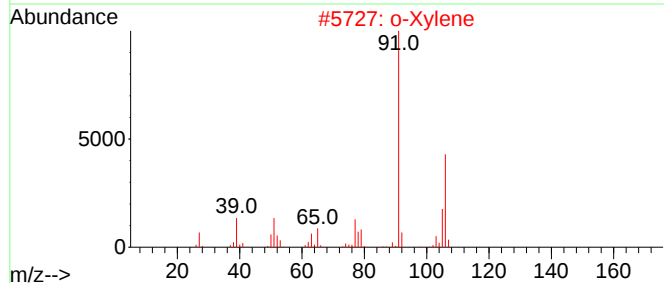
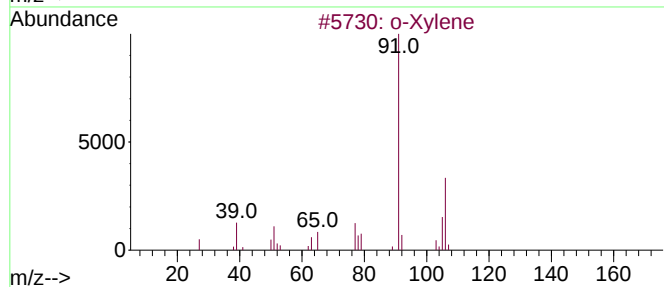
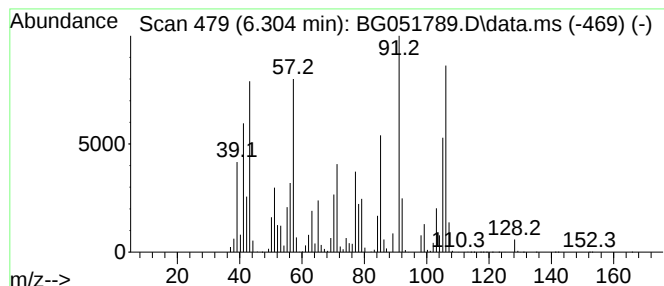
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 o-Xylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.304	55.88 ng/ul	61984900	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	64
2			o-Xylene	106	C8H10	000095-47-6	46
3			p-Xylene	106	C8H10	000106-42-3	46
4			Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	43
5			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

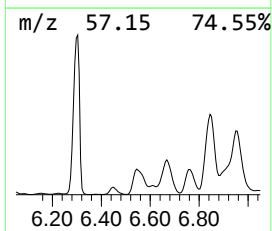
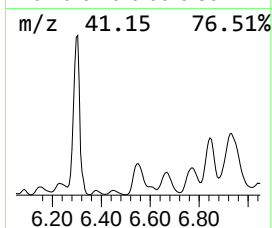
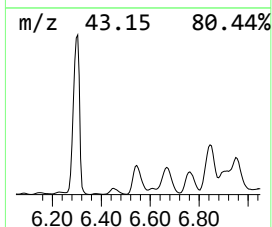
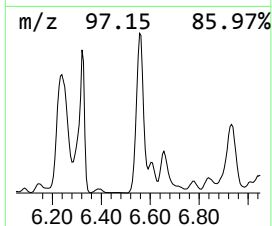
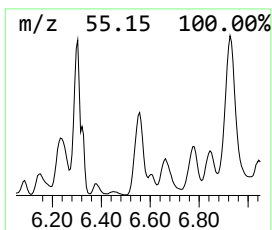
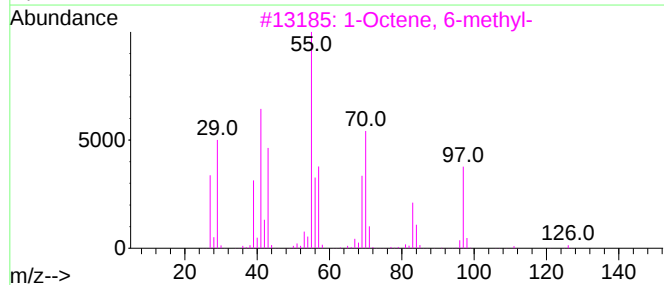
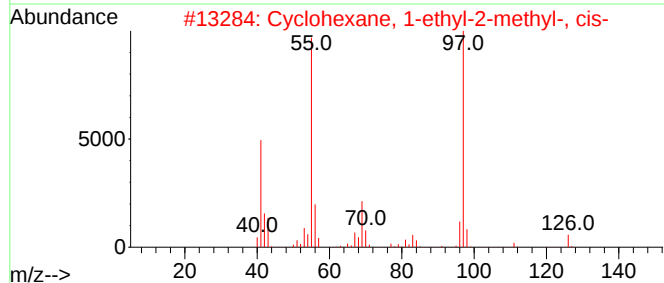
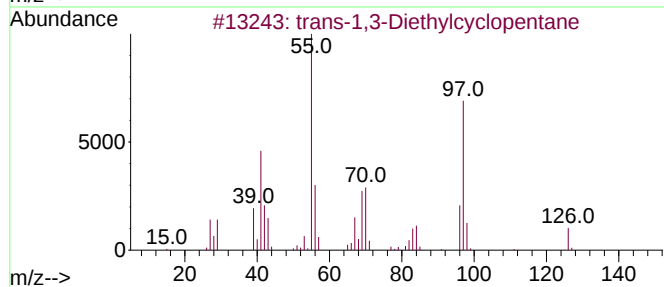
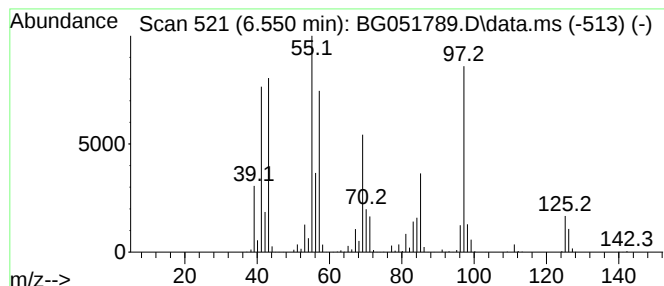
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.550	8.76 ng/ul	9715190	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	53
2		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	53
3		1-Octene, 6-methyl-	126	C9H18	013151-10-5	49
4		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	47
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

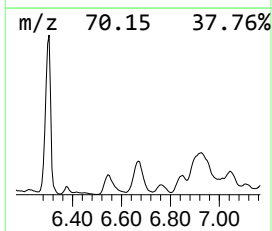
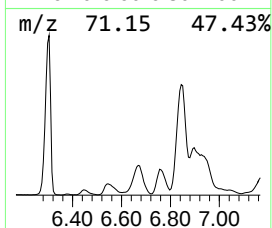
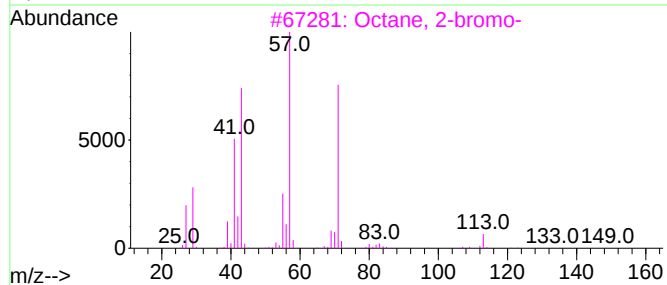
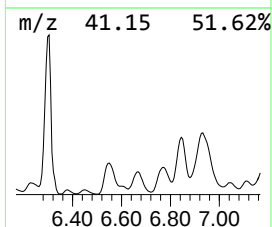
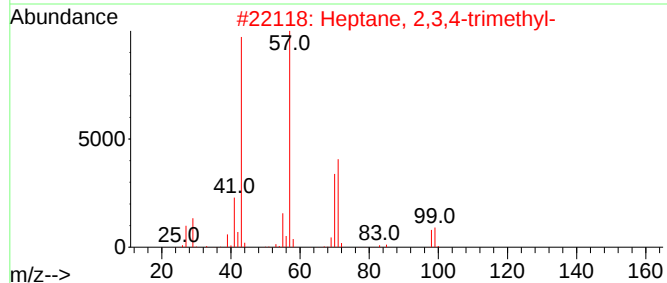
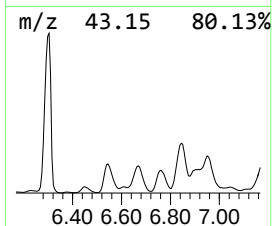
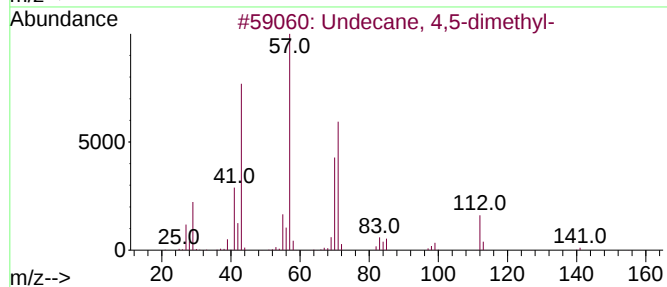
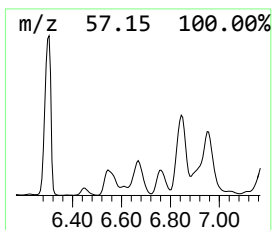
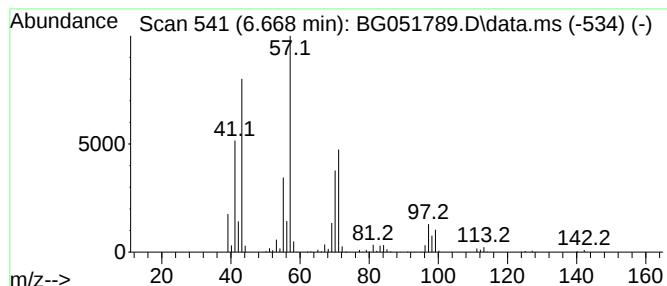
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ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.668	4.23 ng/ul	4690990	1,4-Dichlorobenzene-d4	8.324	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	64
2	Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	64
3	Octane, 2-bromo-	192	C8H17Br	000557-35-7	53
4	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	52
5	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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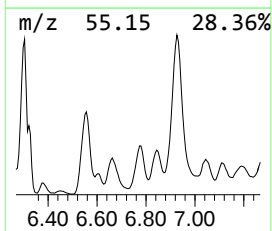
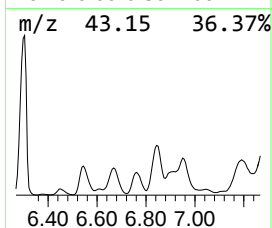
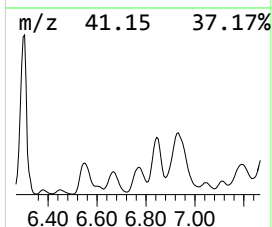
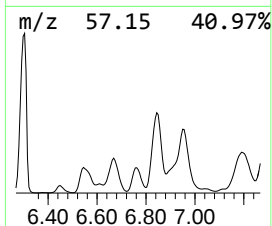
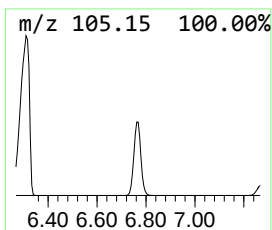
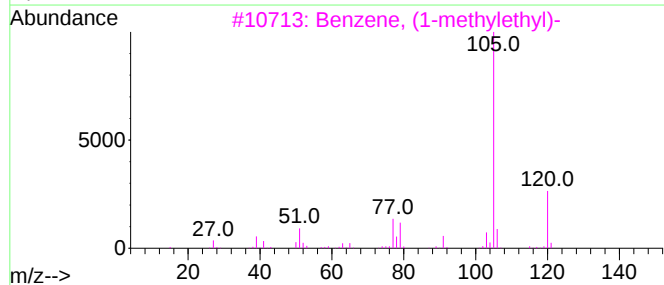
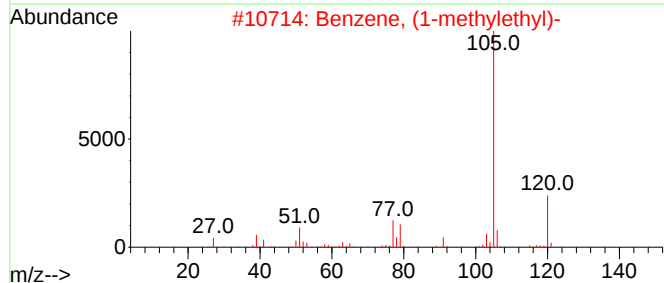
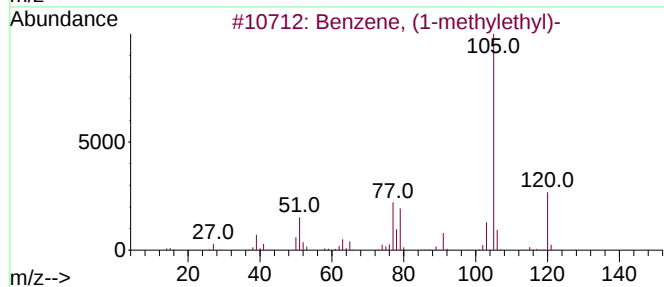
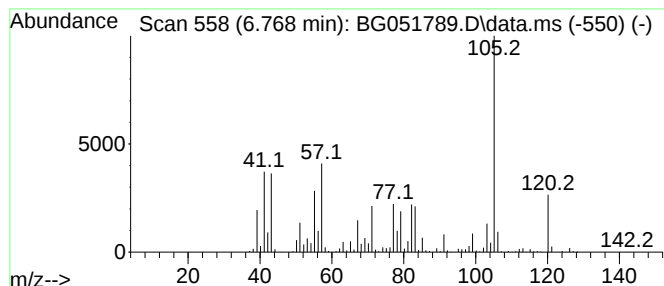
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.768	9.19 ng/ul	10191000	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	95
2			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	55
3			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	55
4			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	50
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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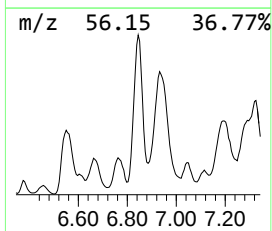
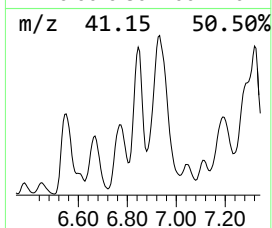
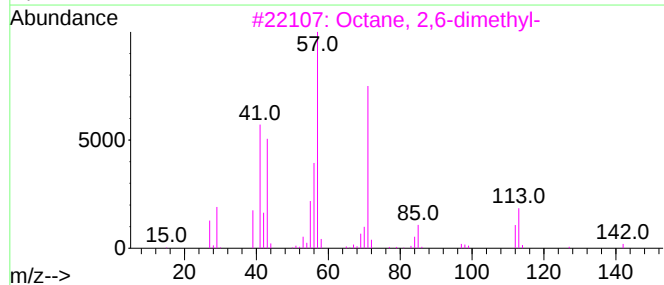
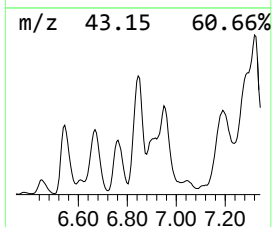
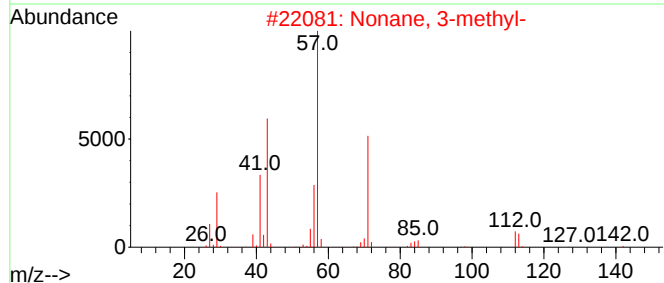
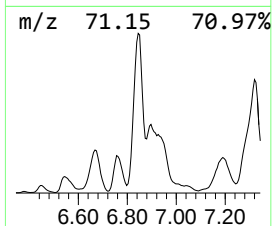
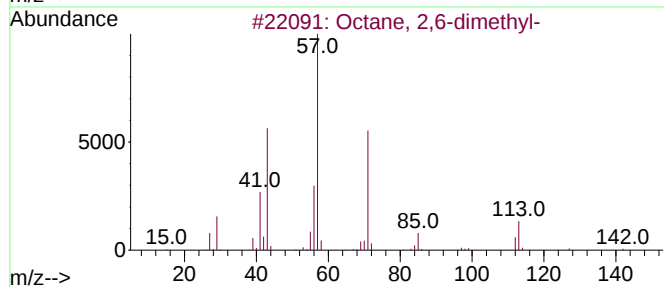
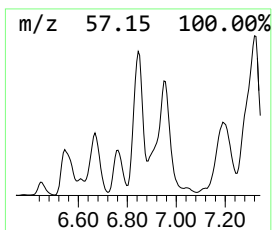
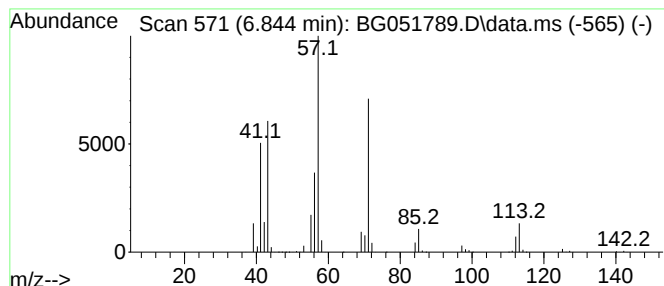
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.844	5.59 ng/ul	6195780	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	87
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	86
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	83
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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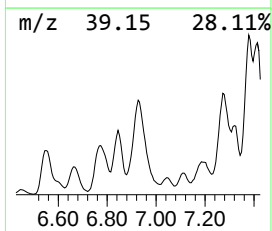
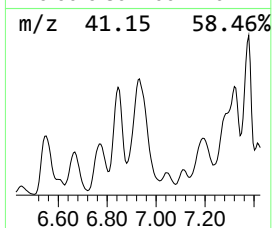
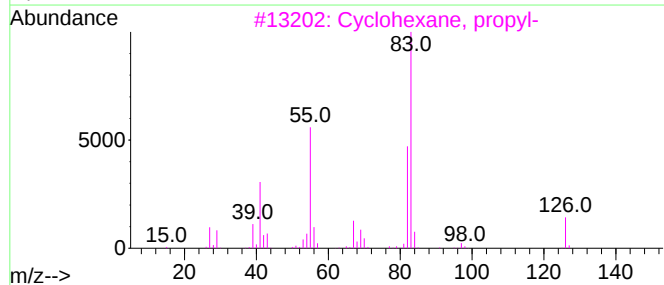
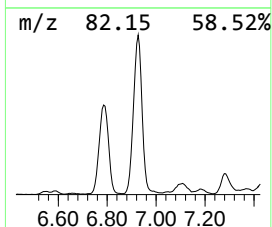
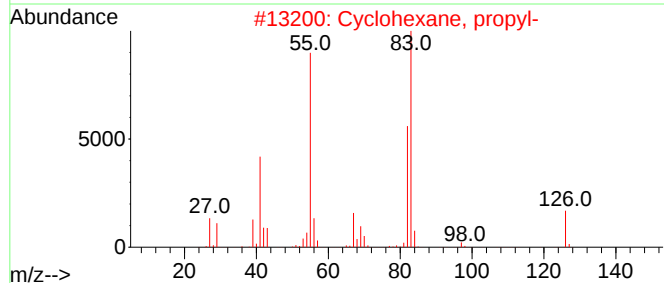
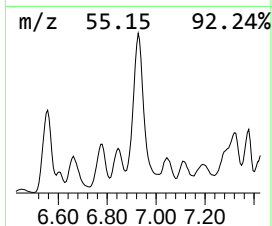
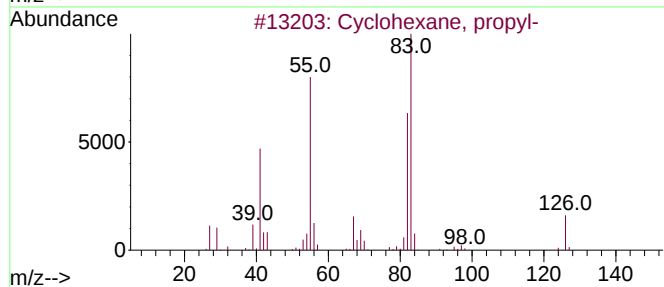
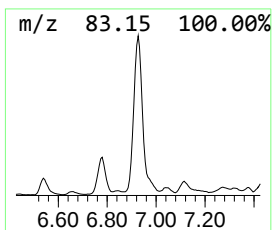
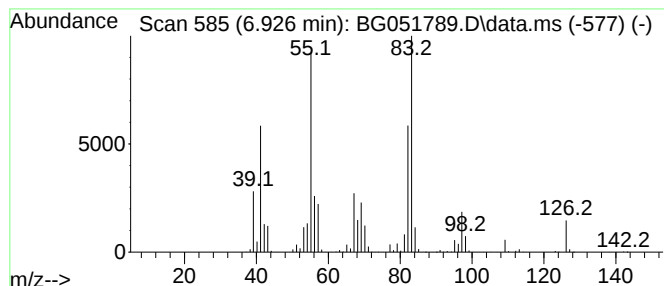
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic6.926 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.926	16.43 ng/ul	18220700	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
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	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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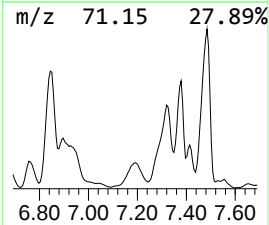
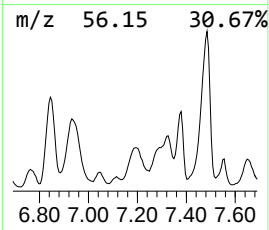
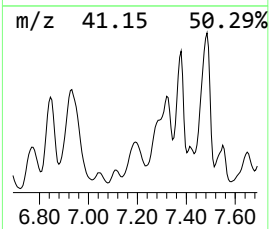
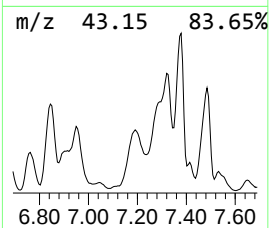
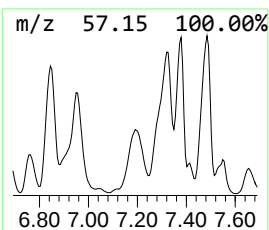
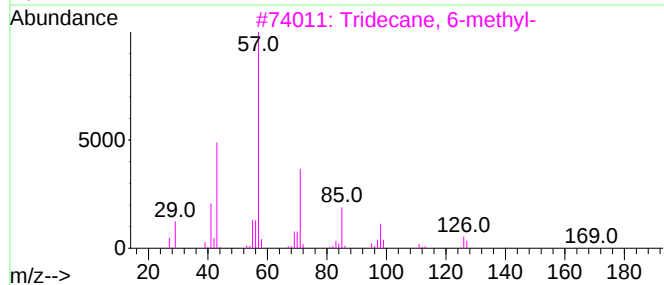
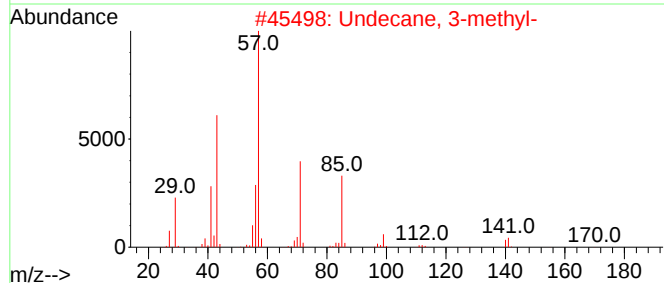
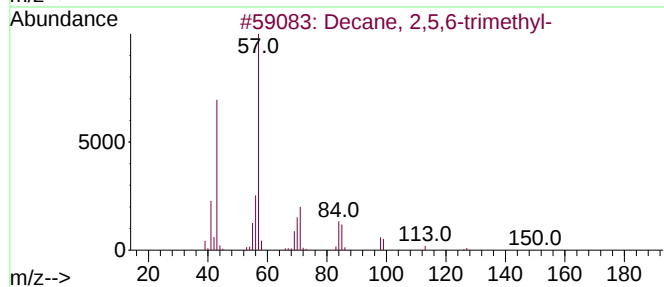
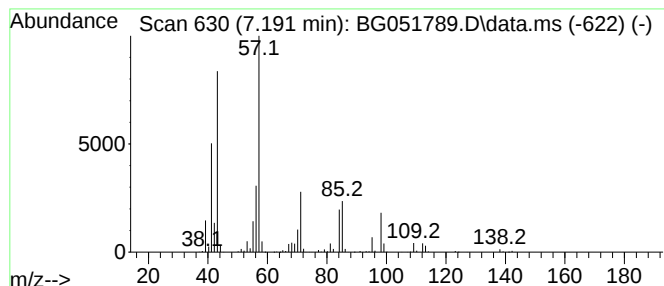
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.191	4.65 ng/ul	5162160	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	80
2			Undecane, 3-methyl-	170	C12H26	001002-43-3	50
3			Tridecane, 6-methyl-	198	C14H30	013287-21-3	47
4			Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1010309-37-4	43
5			Nonane, 2-methyl-	142	C10H22	000871-83-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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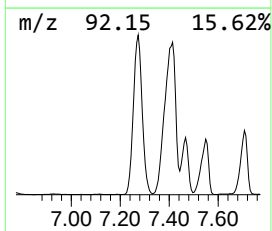
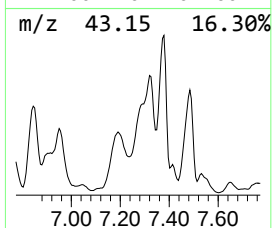
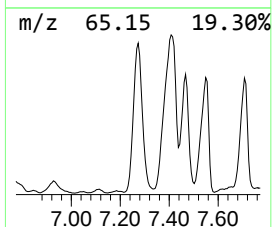
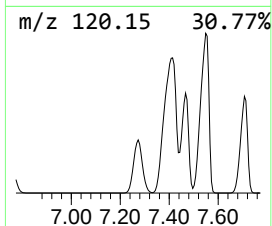
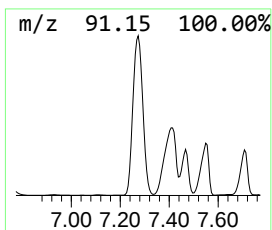
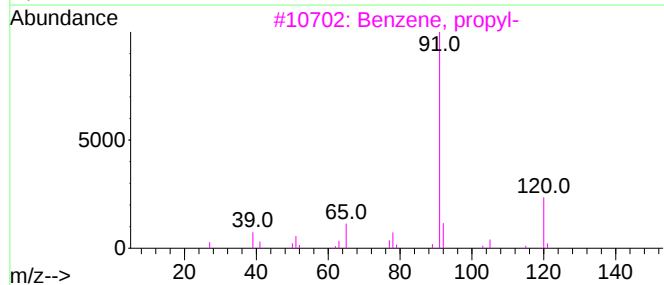
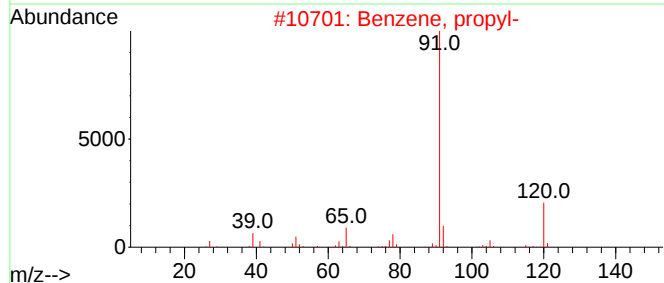
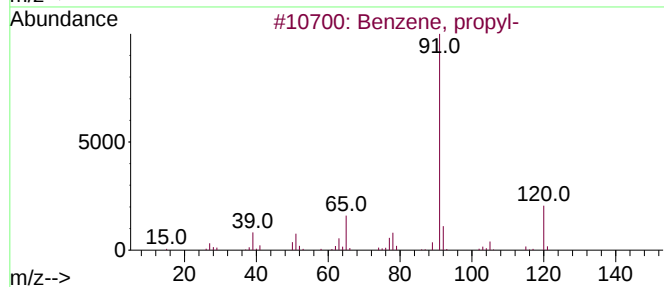
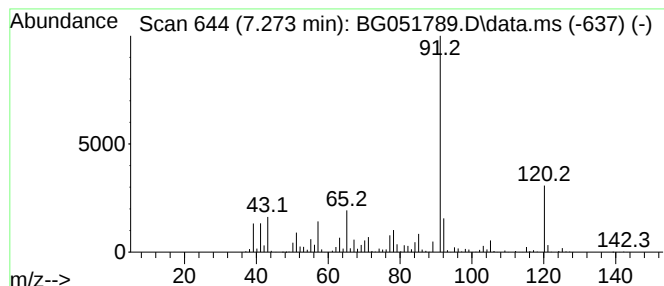
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.273	15.08 ng/ul	16730700	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Benzene, propyl-	120	C9H12	000103-65-1	64
3			Benzene, propyl-	120	C9H12	000103-65-1	64
4			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	59
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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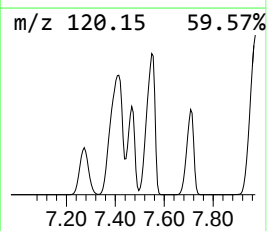
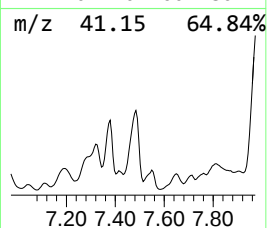
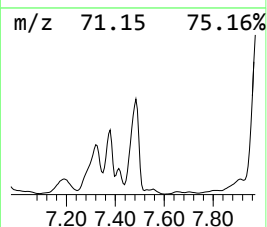
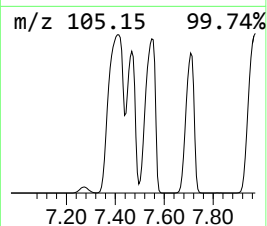
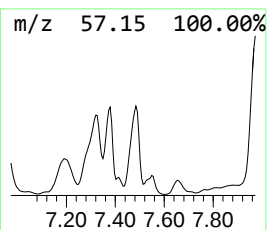
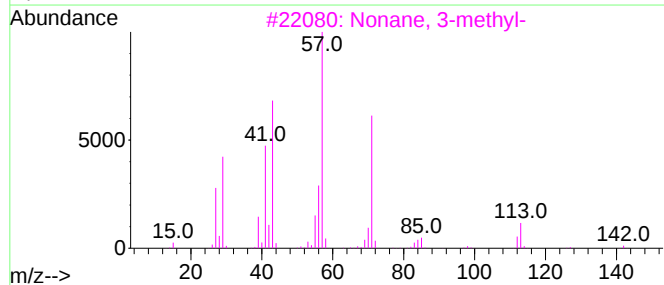
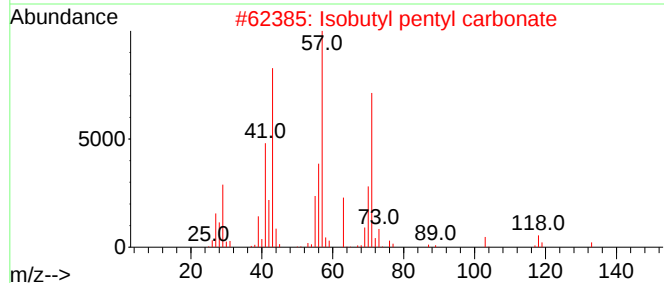
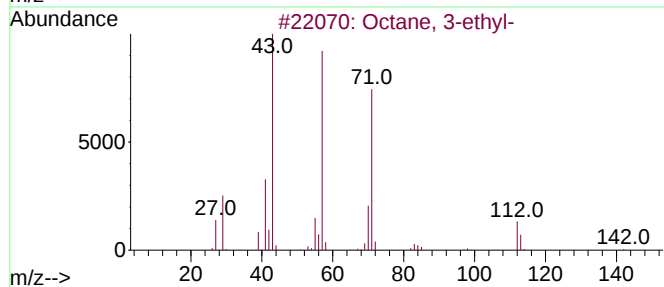
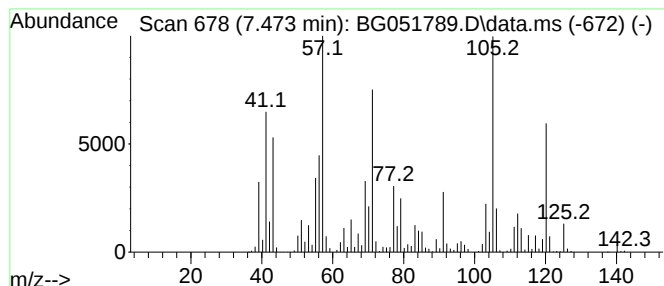
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.473	23.60 ng/ul	26172000	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-			142	C10H22	005881-17-4	38
2	Isobutyl pentyl carbonate			188	C10H20O3	178442-32-5	35
3	Nonane, 3-methyl-			142	C10H22	005911-04-6	35
4	Benzene, 1-ethyl-2-methyl-			120	C9H12	000611-14-3	35
5	Benzene, 1-ethyl-4-methyl-			120	C9H12	000622-96-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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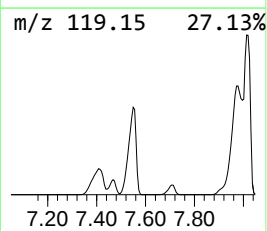
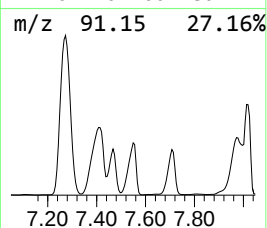
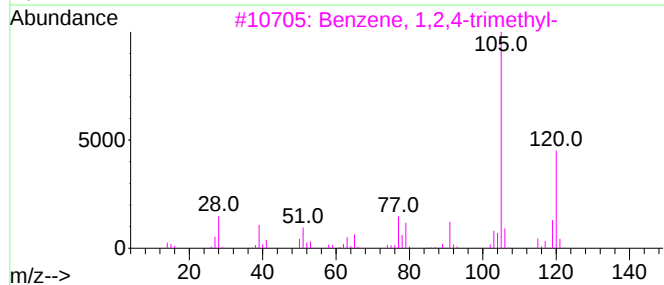
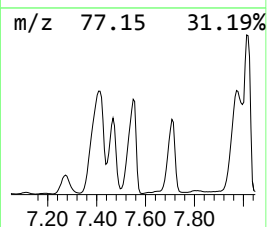
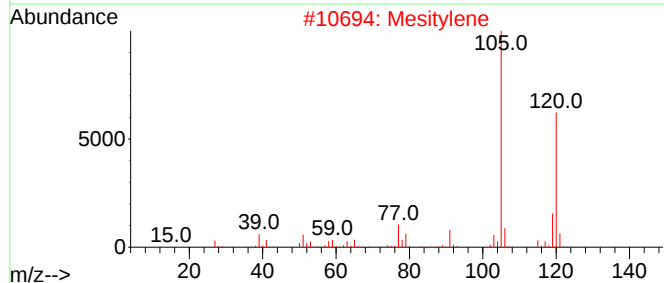
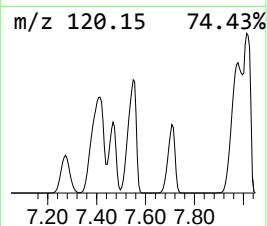
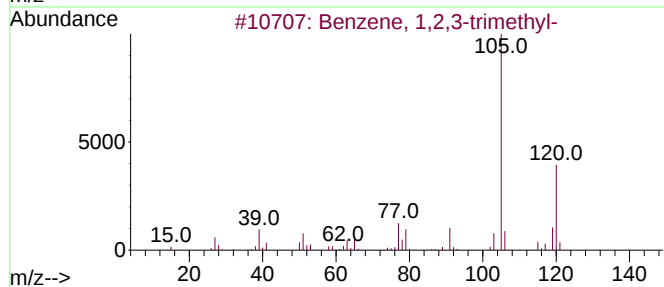
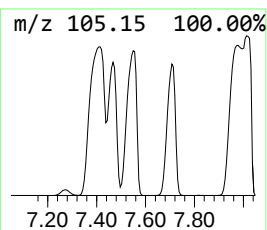
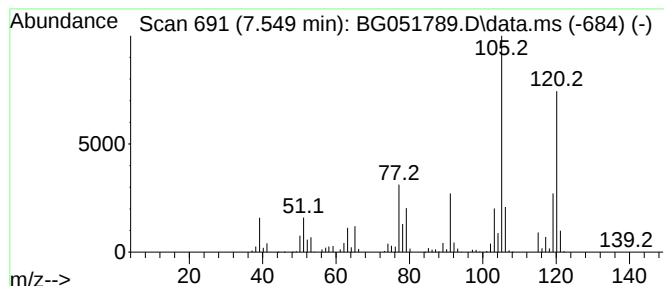
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.549	22.42 ng/ul	24866900	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
2			Mesitylene	120	C9H12	000108-67-8	87
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	81
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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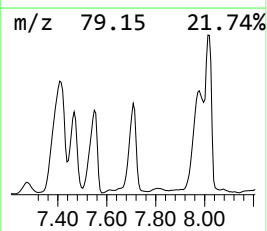
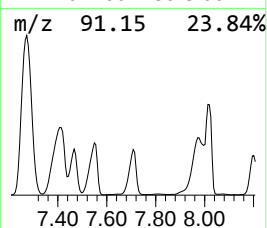
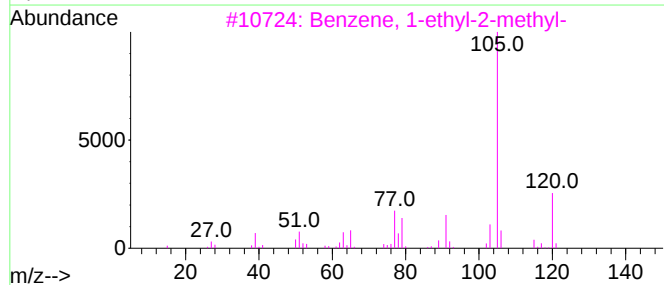
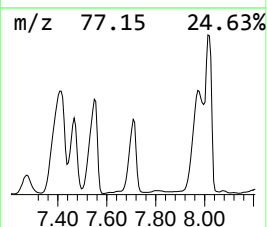
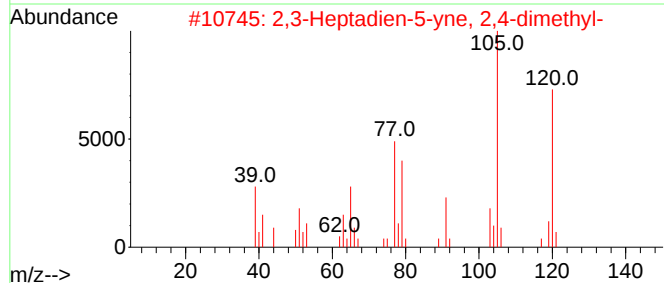
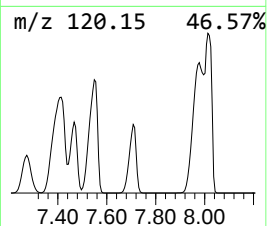
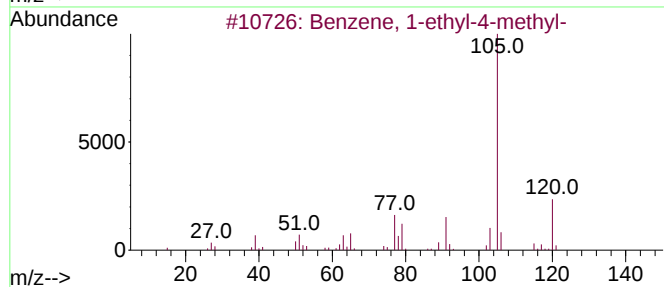
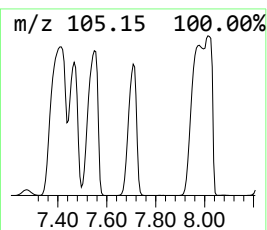
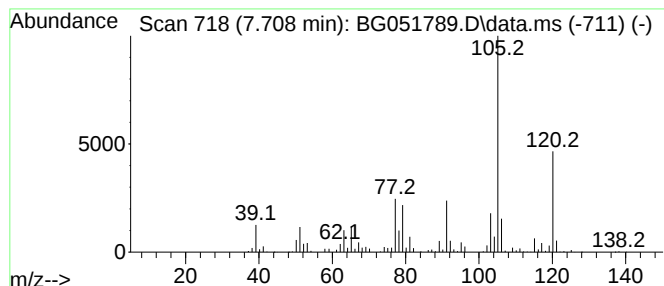
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzene, 1-ethyl-4-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.708	13.20 ng/ul	14635600	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2			2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	90
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	89
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	86
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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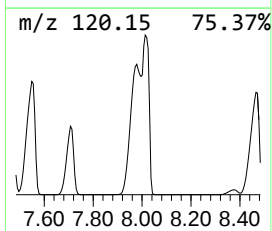
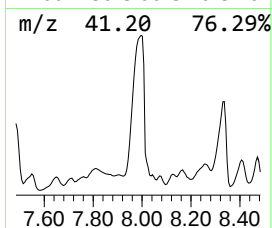
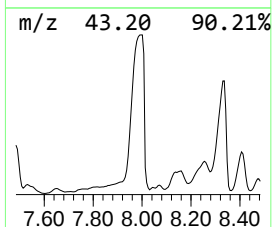
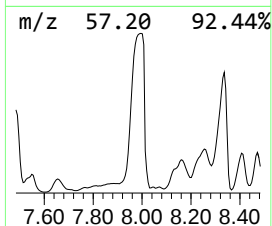
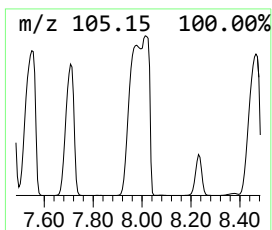
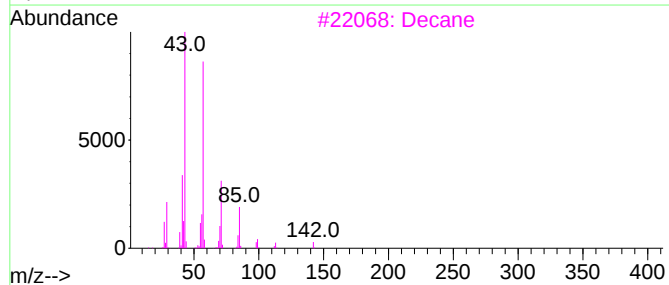
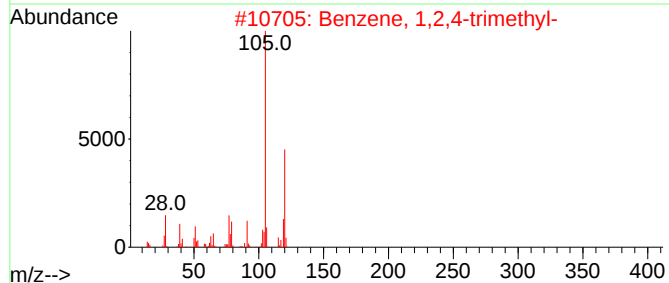
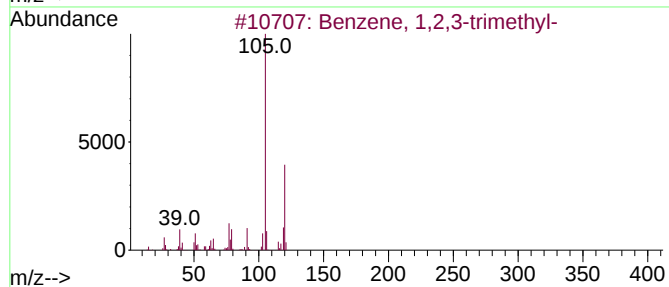
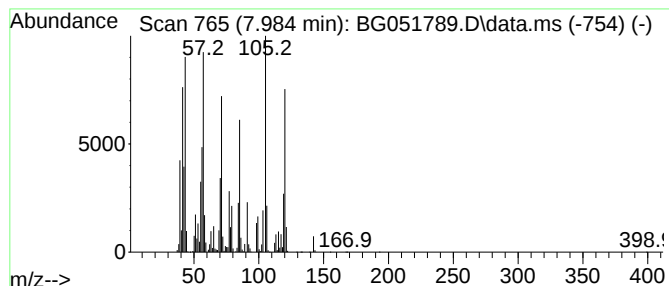
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.984	59.49 ng/ul	65985100	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	83
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	62
3			Decane	142	C10H22	000124-18-5	43
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	42
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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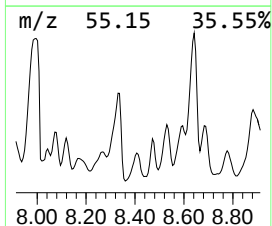
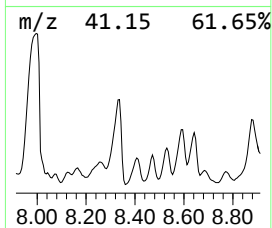
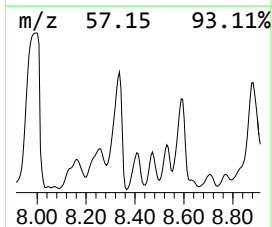
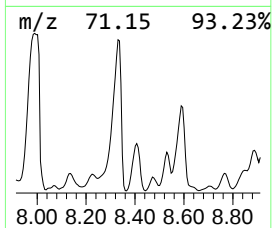
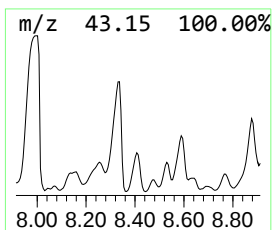
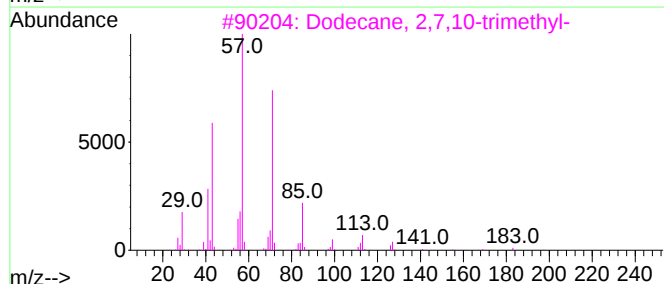
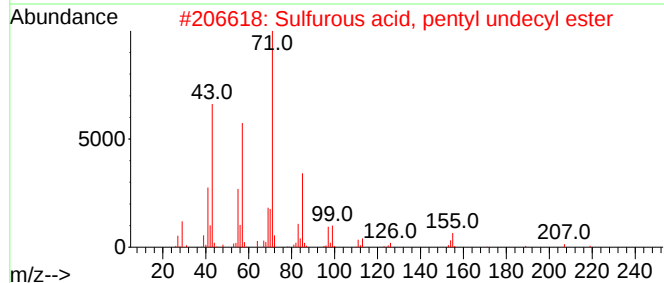
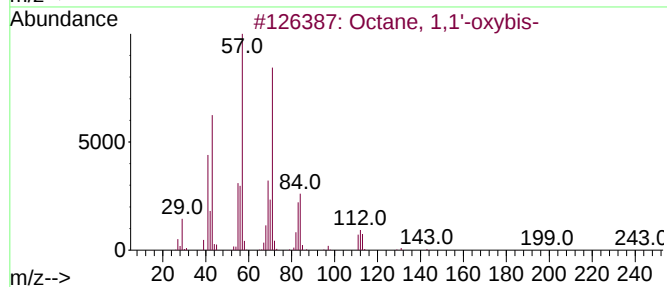
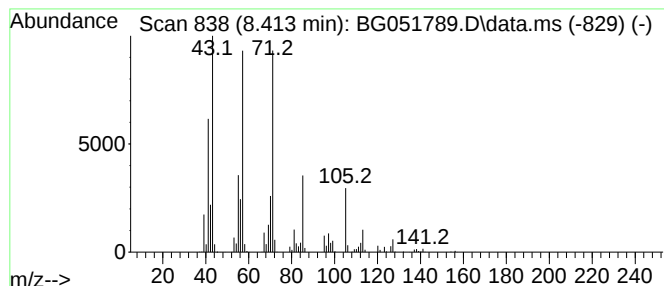
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Octane, 1,1'-oxybis- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.413	6.21 ng/ul	6884660	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	64
2			Sulfurous acid, pentyl undecyl e...	306	C16H34O3S	1000309-14-4	59
3			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	58
4			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	58
5			Hexadecyl pentyl ether	312	C21H44O	1000406-41-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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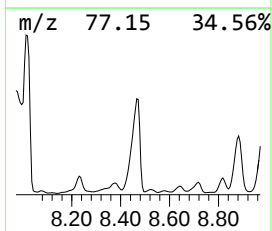
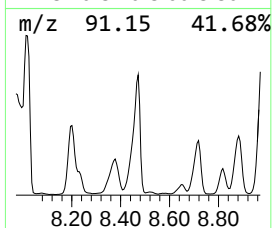
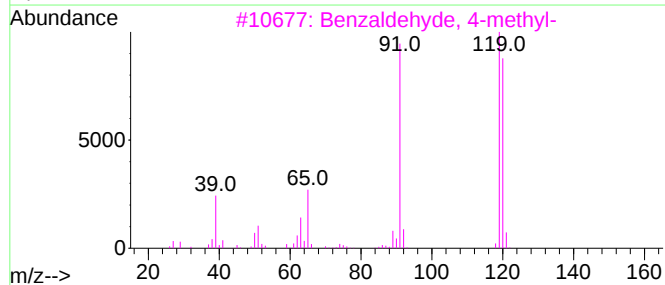
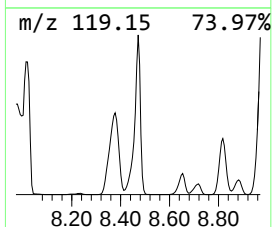
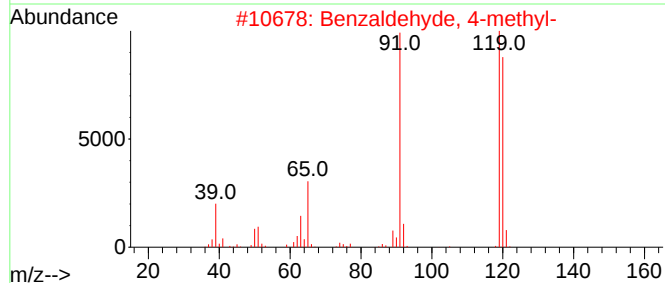
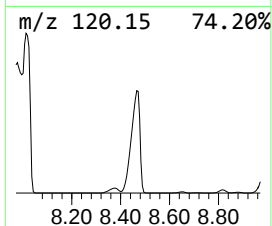
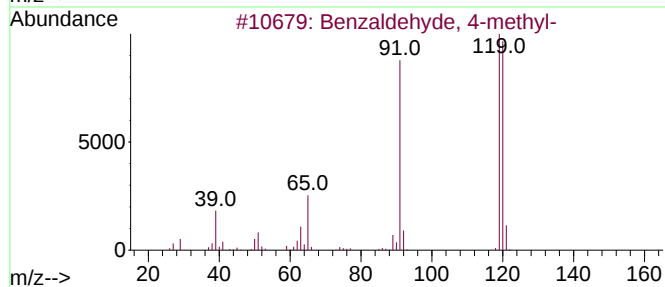
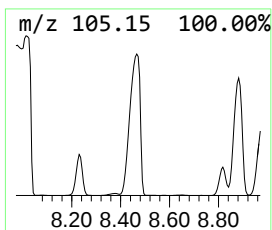
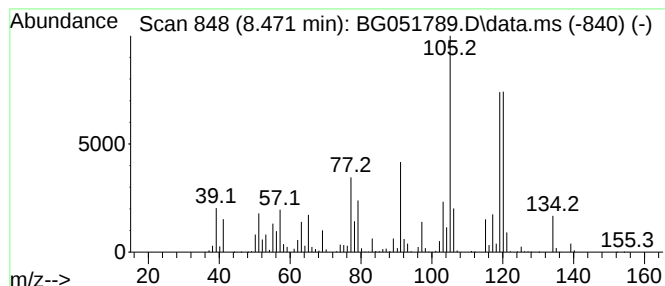
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzaldehyde, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.471	23.03 ng/ul	25546700	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	78
2			Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	60
3			Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	53
4			2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	49
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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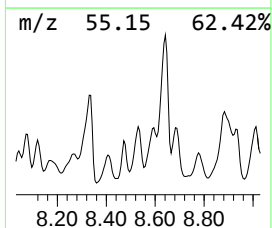
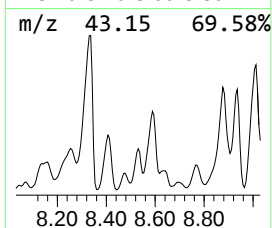
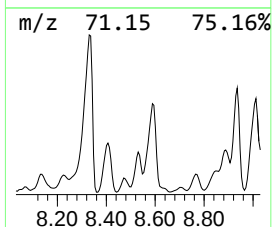
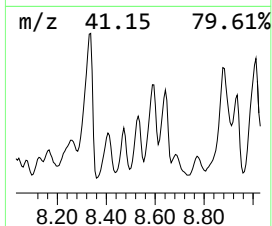
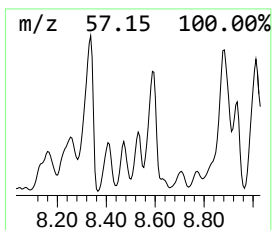
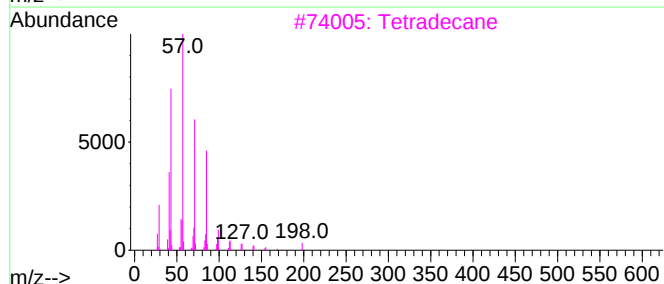
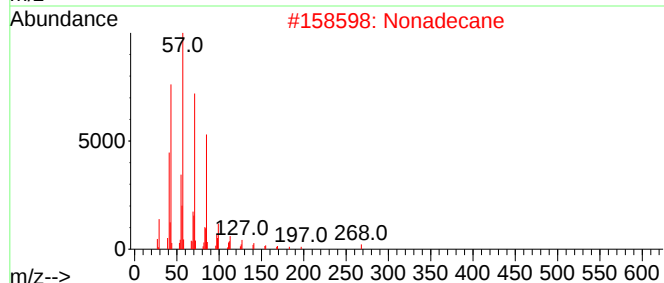
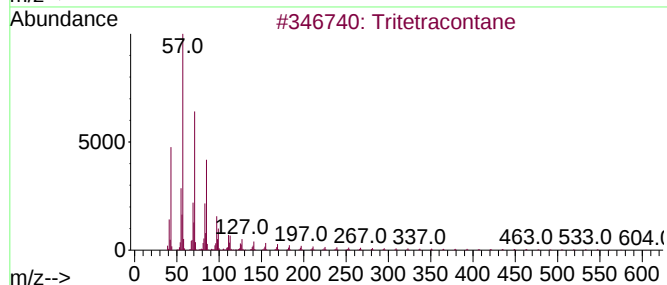
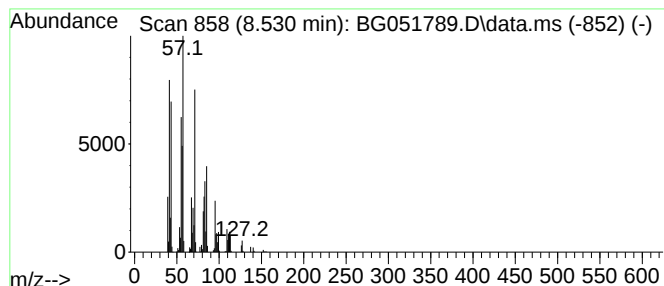
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.530	5.73 ng/ul	6359030	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tritetracontane	605	C43H88	007098-21-7	50
2		Nonadecane	268	C19H40	000629-92-5	49
3		Tetradecane	198	C14H30	000629-59-4	49
4		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	46
5		Pentadecane	212	C15H32	000629-62-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

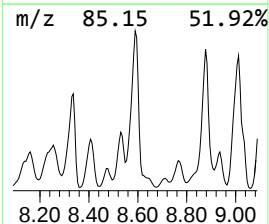
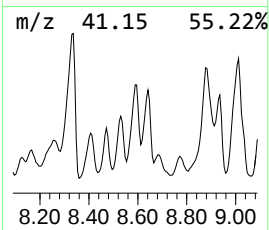
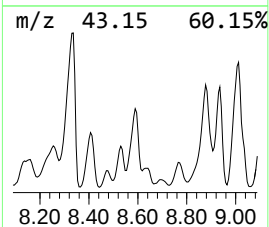
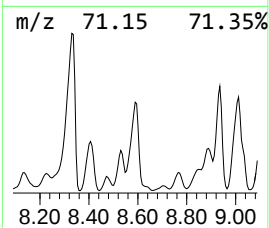
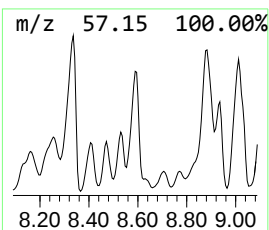
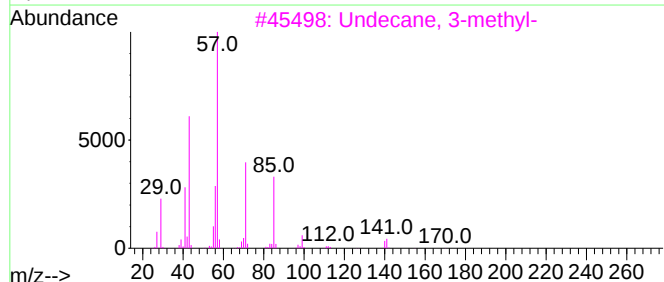
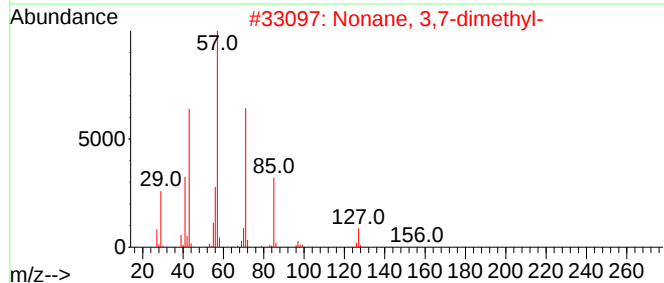
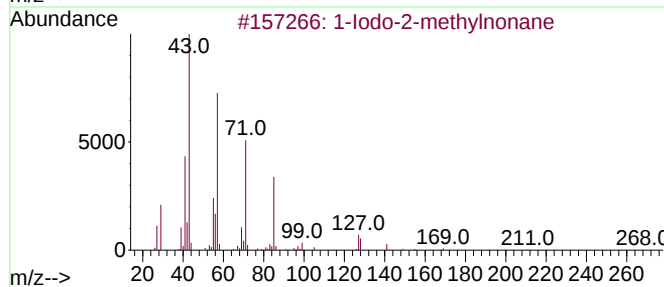
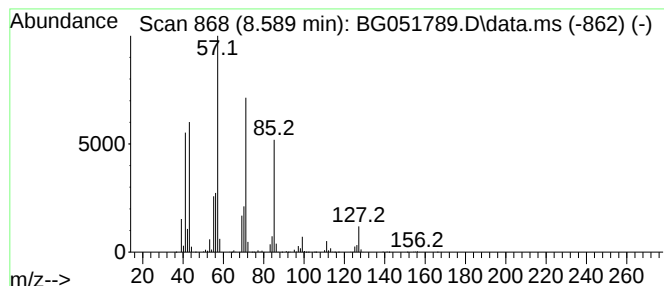
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1-Iodo-2-methylnonane Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.589	9.51 ng/ul	10546400	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	72
2		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	70
3		Undecane, 3-methyl-	170	C12H26	001002-43-3	64
4		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	58
5		Octane	114	C8H18	000111-65-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
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Instrument :
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ClientSampleId :
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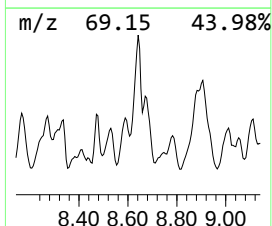
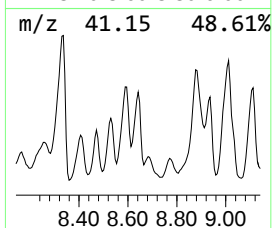
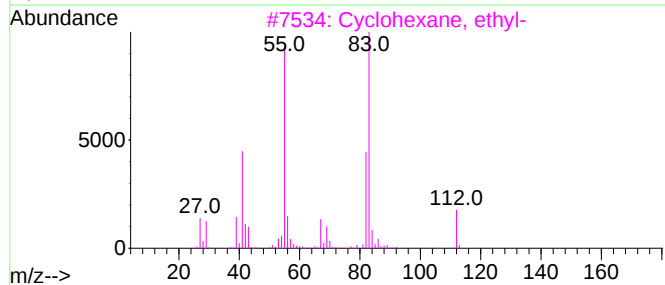
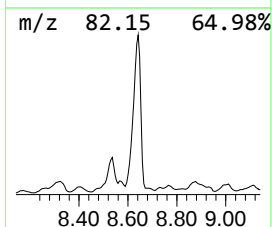
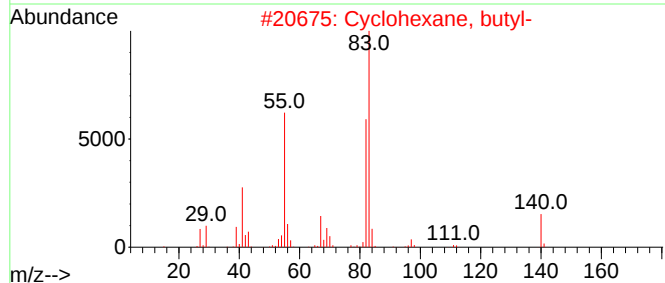
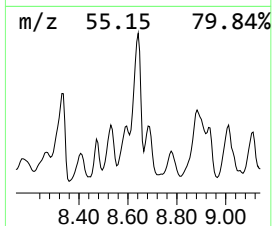
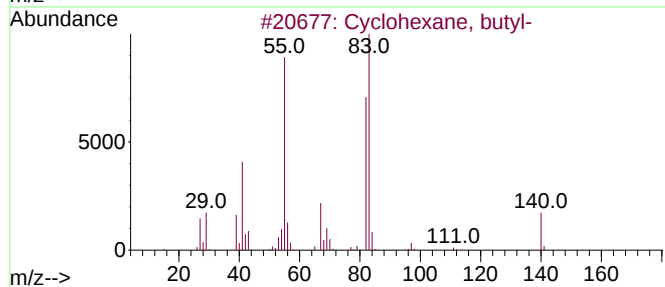
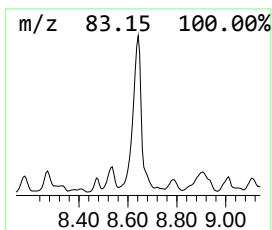
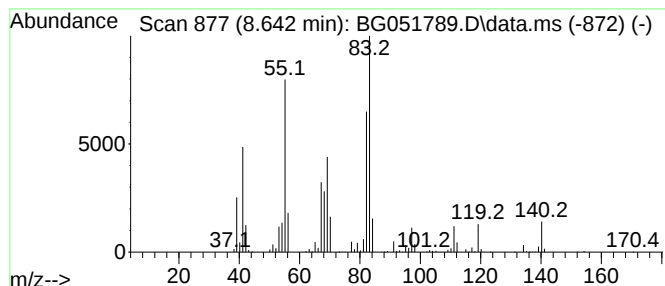
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Cyclic8.642 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.642	8.82 ng/ul	9779200	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, butyl-	140	C10H20	001678-93-9	64
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	59
3		Cyclohexane, ethyl-	112	C8H16	001678-91-7	53
4		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53
5		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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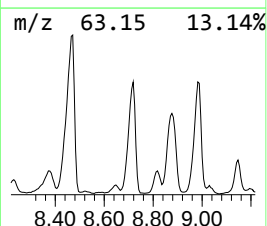
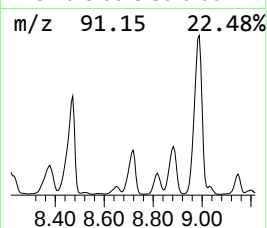
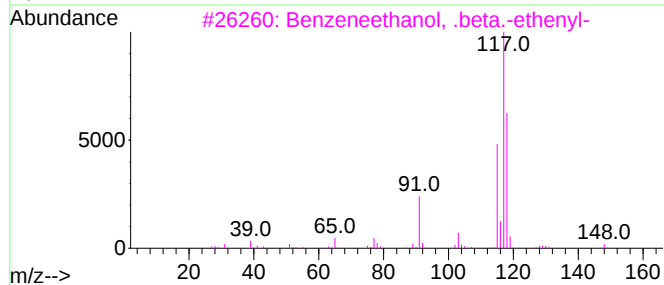
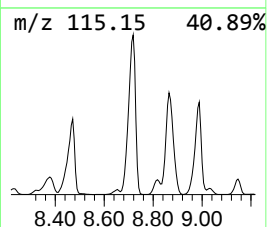
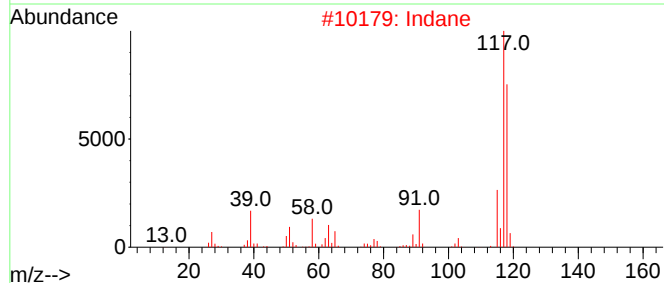
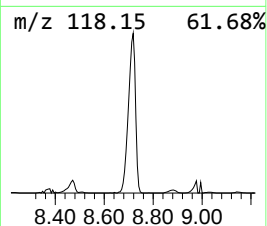
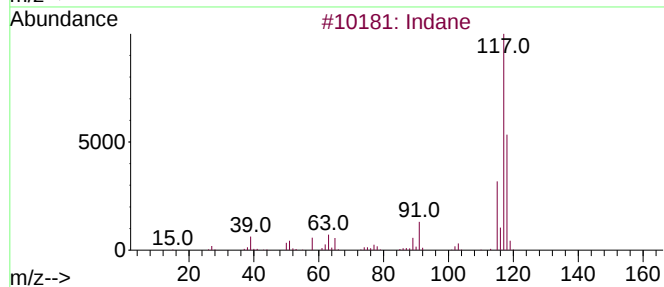
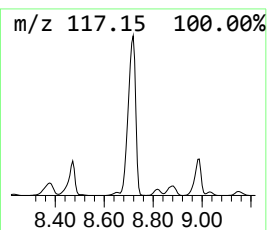
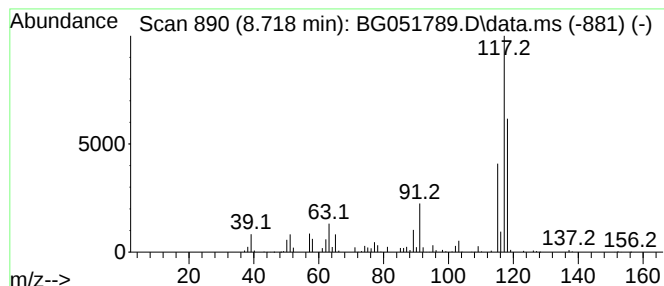
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Indane Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.718	10.63 ng/ul	11791500	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	76
2	Indane			118	C9H10	000496-11-7	76
3	Benzeneethanol, .beta.-ethenyl-			148	C10H12O	006052-63-7	72
4	Benzene, (2-bromocyclopropyl)-			196	C9H9Br	036617-02-4	64
5	Benzene, 2-propenyl-			118	C9H10	000300-57-2	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
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Instrument :
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ClientSampleId :
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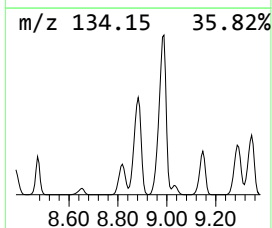
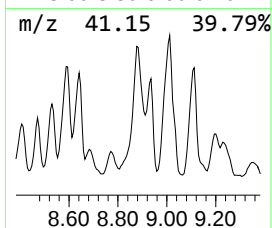
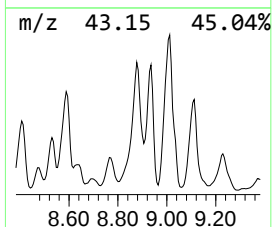
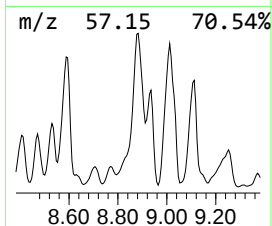
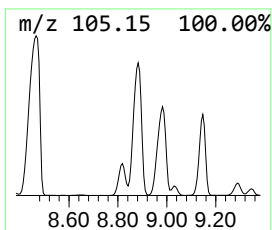
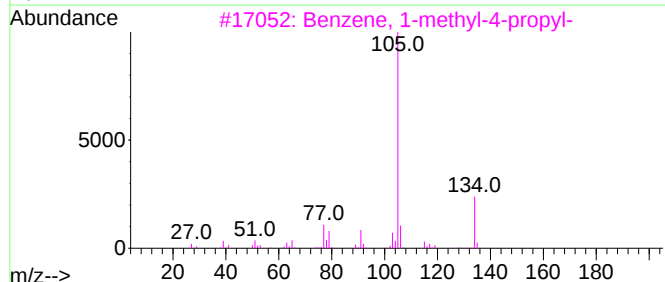
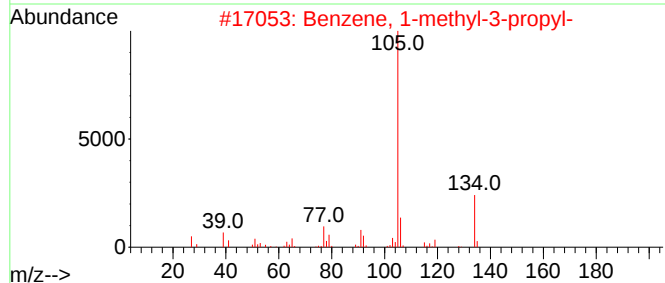
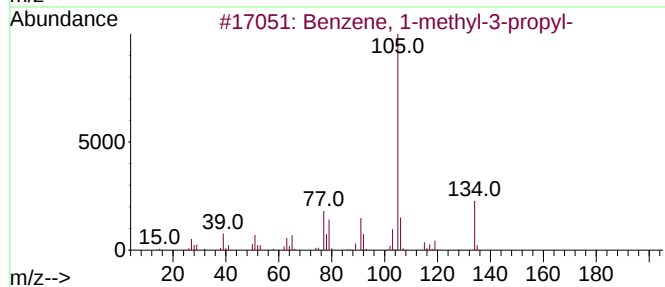
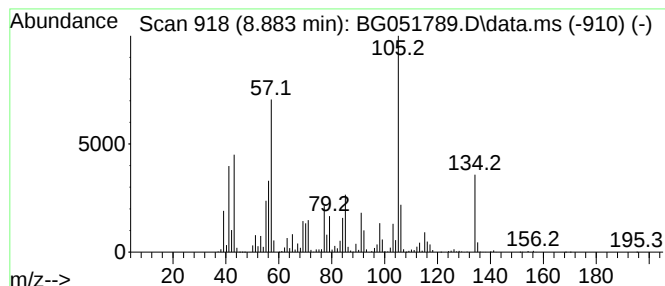
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 1-methyl-3-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.883	24.93 ng/ul	27654100	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	55
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	50
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	46
4			Benzeneacetaldehyde, 4-methyl-	134	C9H10O	000104-09-6	45
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Sample : M5215-04
Misc :
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Instrument :
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ClientSampleId :
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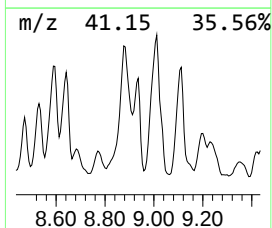
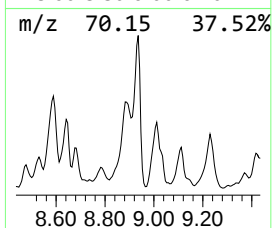
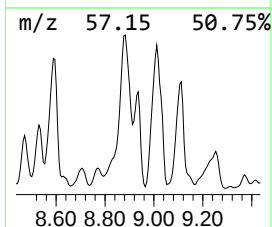
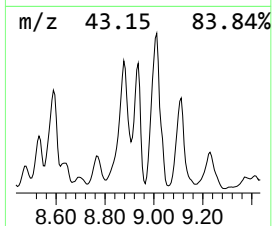
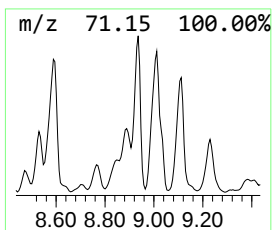
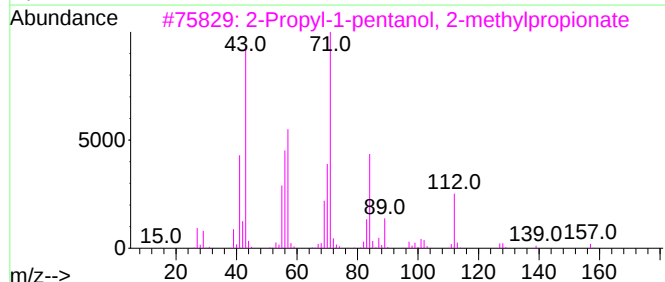
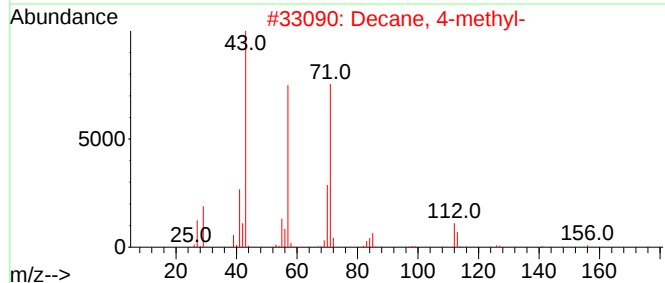
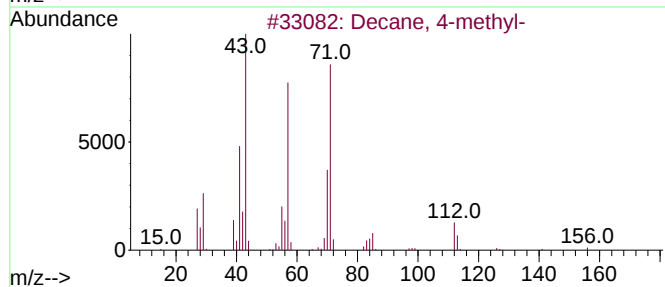
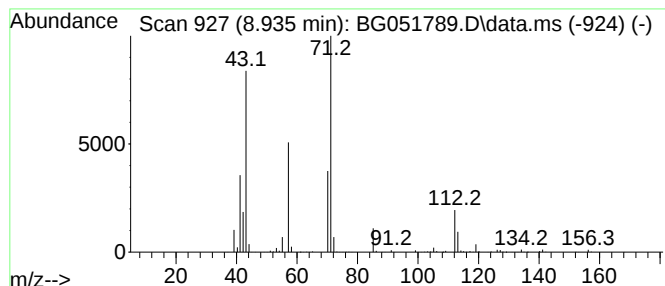
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.935	4.99 ng/ul	5530300	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 4-methyl-	156	C11H24	002847-72-5	80
2		Decane, 4-methyl-	156	C11H24	002847-72-5	74
3		2-Propyl-1-pentanol, 2-methylpro...	200	C12H24O2	1000447-36-4	72
4		Octane, 3-ethyl-	142	C10H22	005881-17-4	64
5		Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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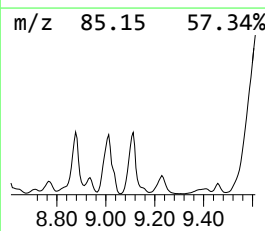
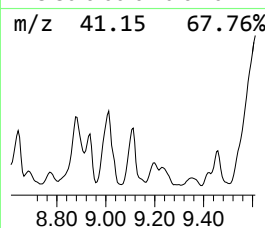
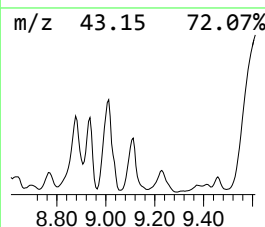
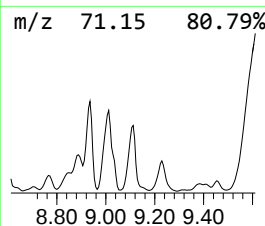
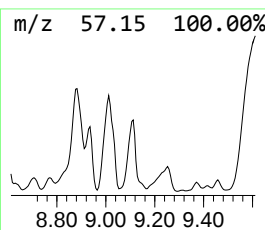
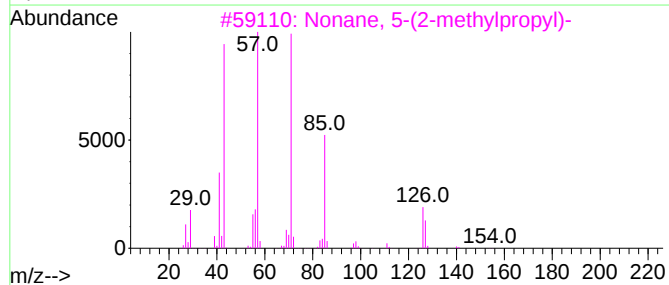
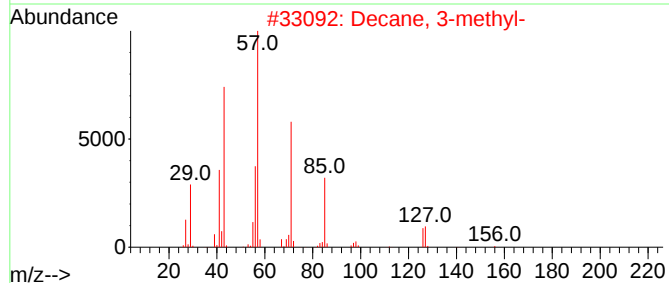
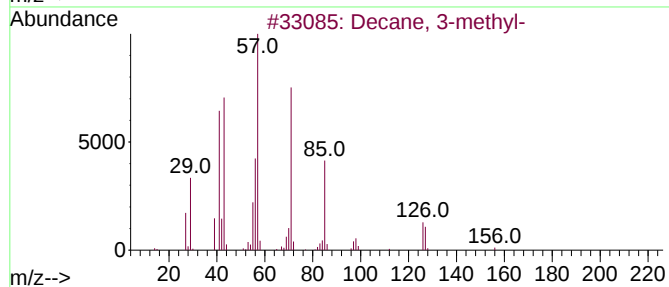
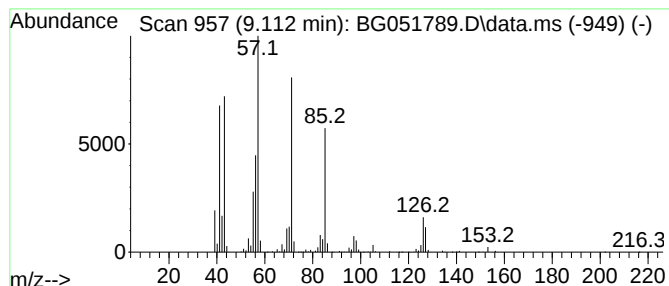
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.112	10.49 ng/ul	11639300	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	90
2			Decane, 3-methyl-	156	C11H24	013151-34-3	87
3			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	72
4			Nonane, 5-butyl-	184	C13H28	017312-63-9	72
5			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

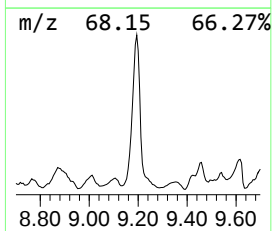
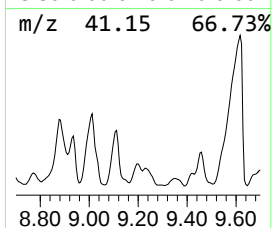
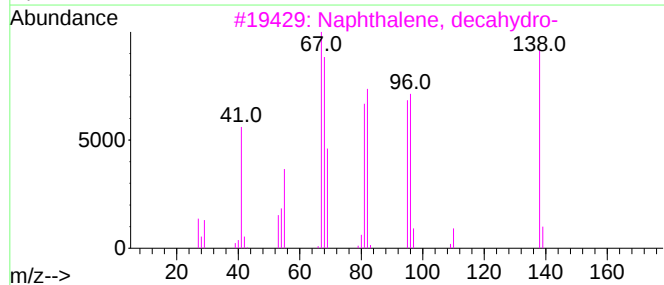
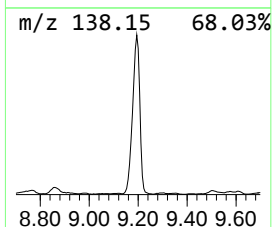
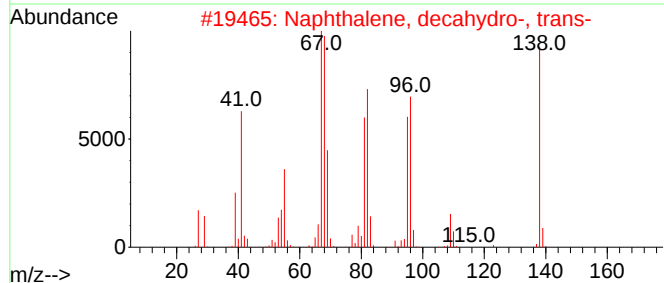
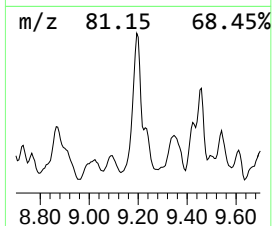
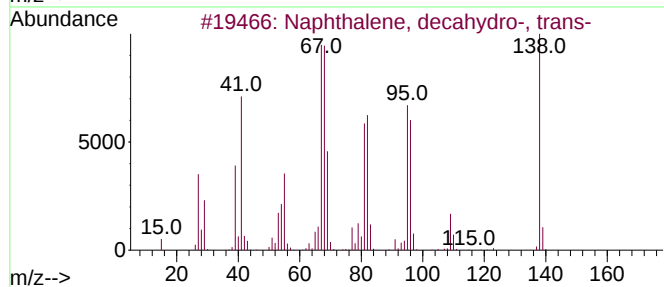
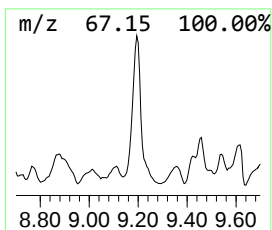
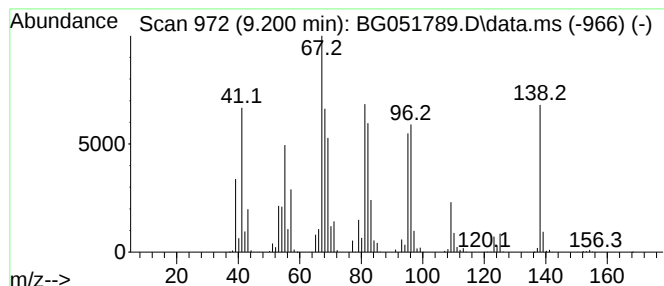
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Naphthalene, decahydro-, tr... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.200	4.79 ng/ul	5315660	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
3			Naphthalene, decahydro-	138	C10H18	000091-17-8	95
4			Naphthalene, decahydro-	138	C10H18	000091-17-8	95
5			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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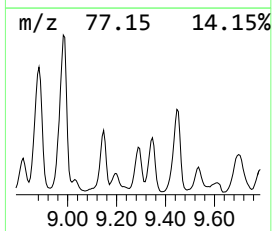
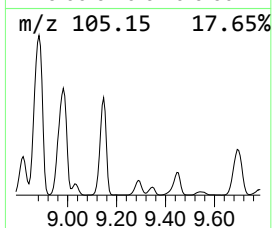
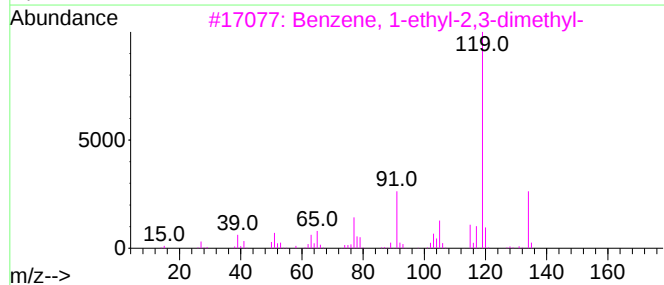
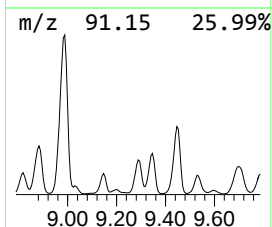
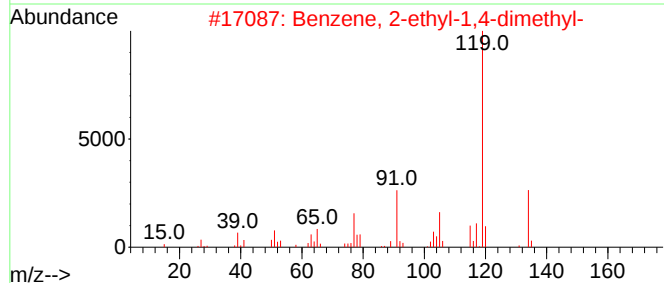
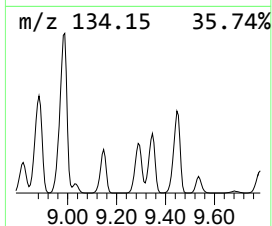
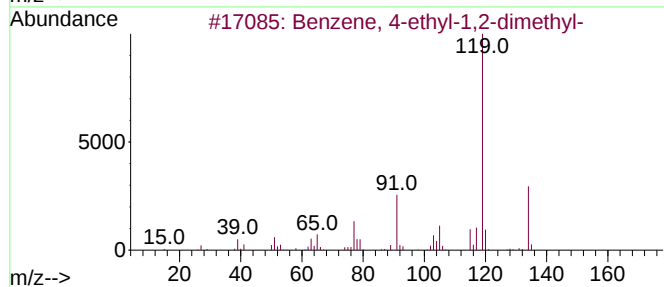
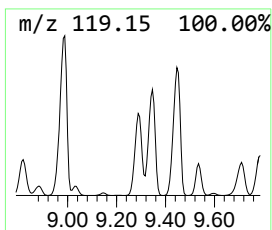
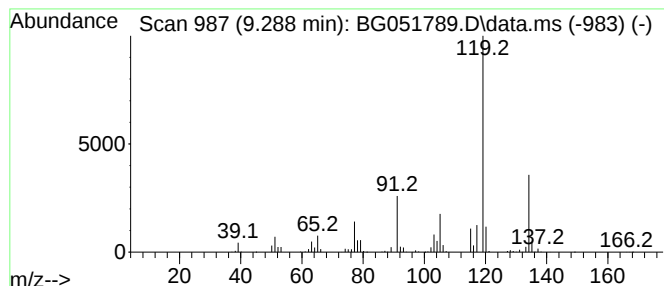
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.288	3.05 ng/ul	3385380	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	97
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

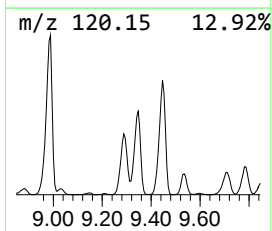
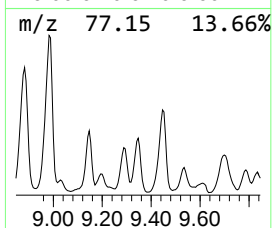
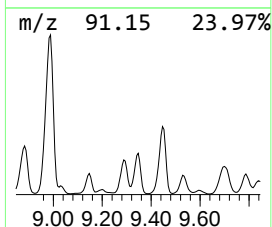
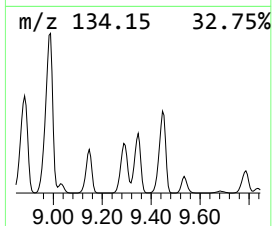
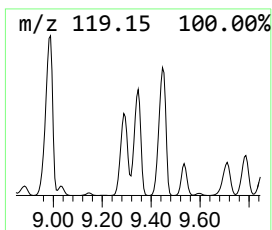
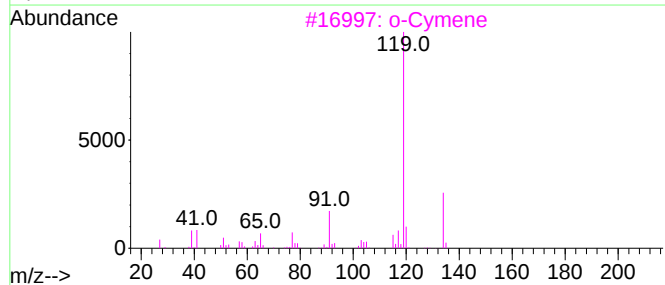
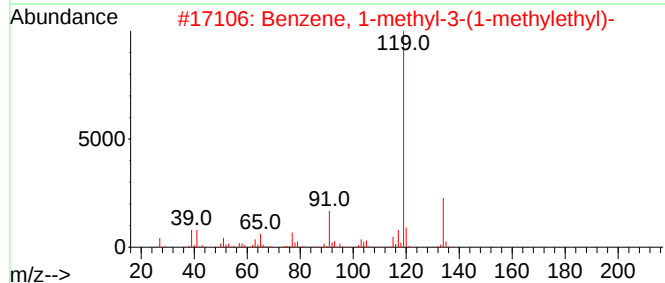
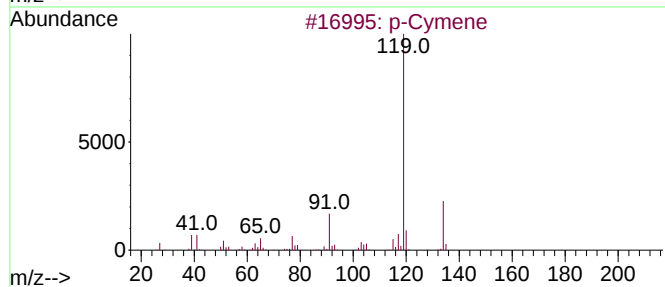
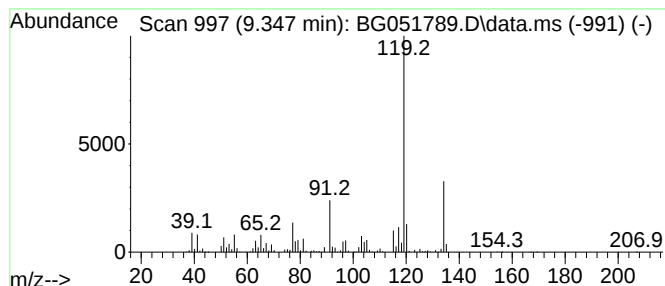
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 p-Cymene Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.347	7.90 ng/ul	8759410	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

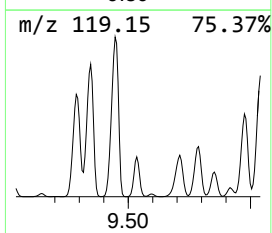
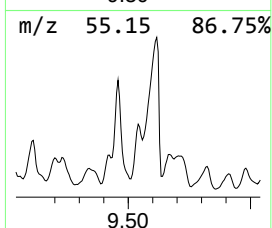
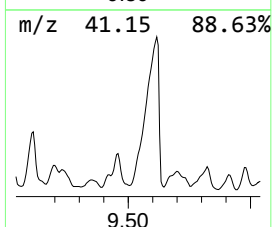
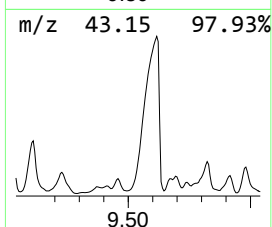
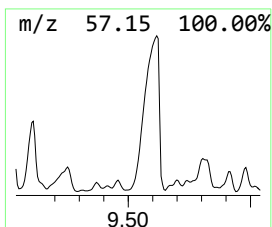
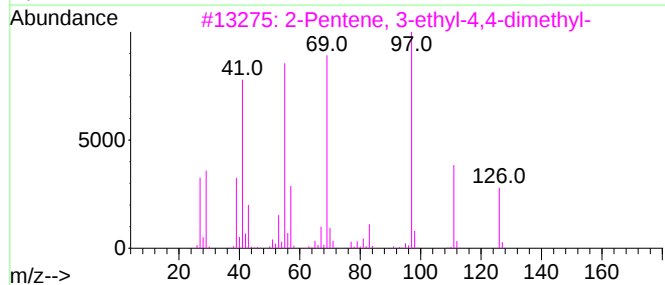
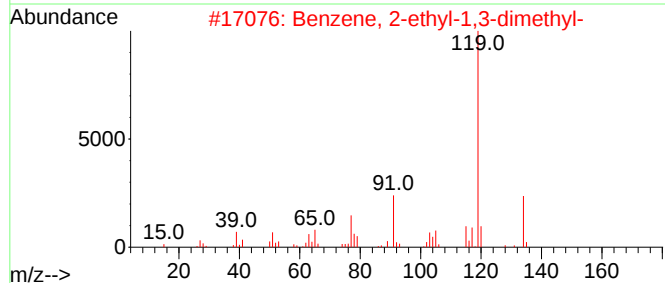
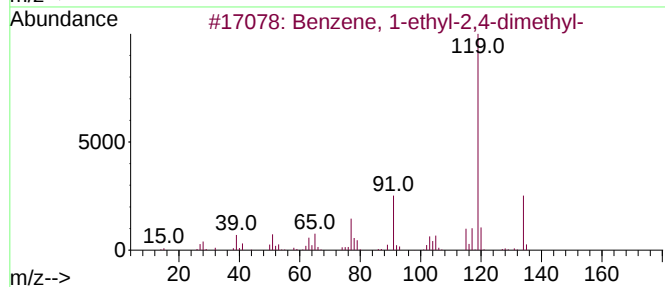
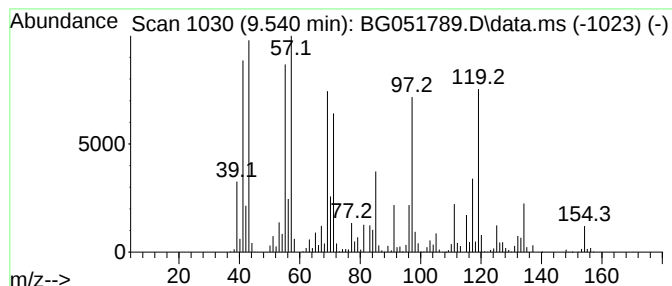
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.540	7.37 ng/ul	8171540	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60
3			2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	42
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	35
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

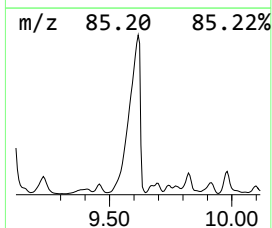
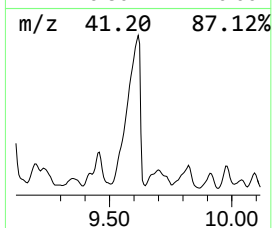
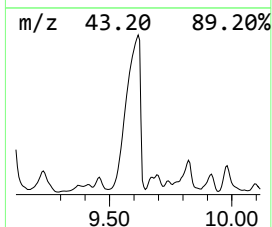
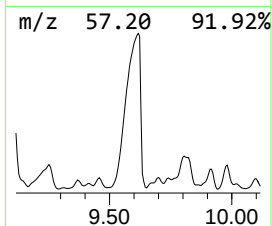
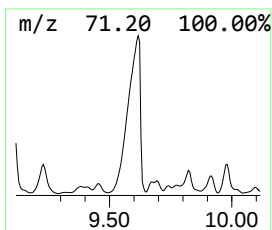
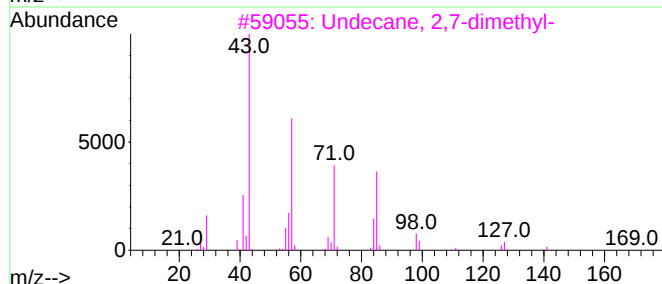
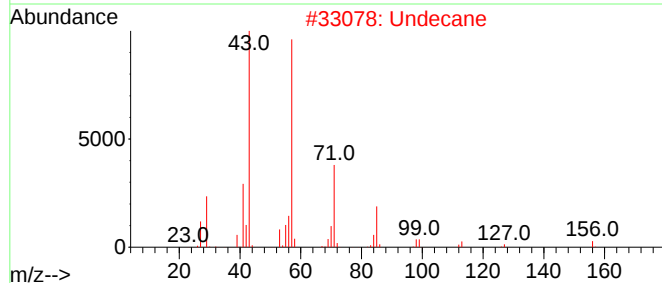
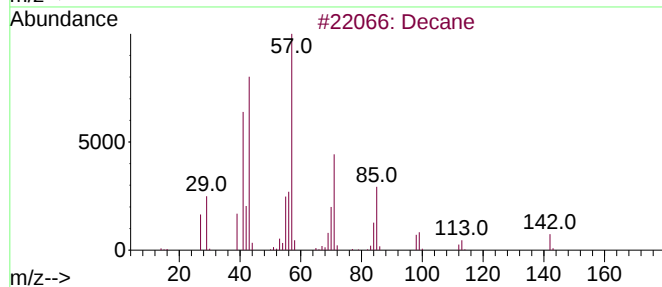
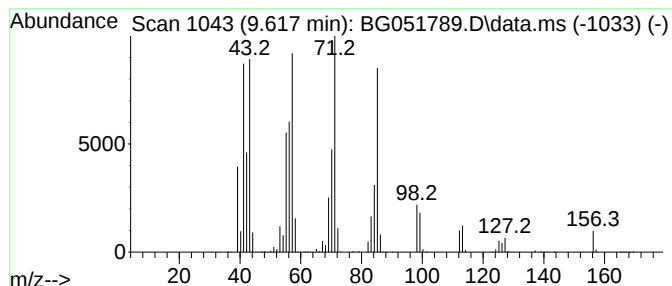
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.617	46.85 ng/ul	51968400	1,4-Dichlorobenzene-d4	8.324

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane	142	C10H22	000124-18-5	90
2		Undecane	156	C11H24	001120-21-4	58
3		Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	52
4		Octane	114	C8H18	000111-65-9	43
5		1-Octene	112	C8H16	000111-66-0	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

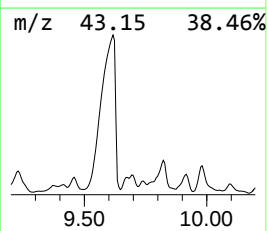
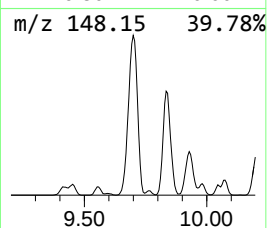
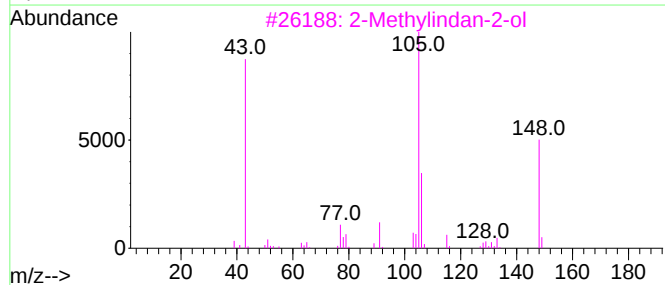
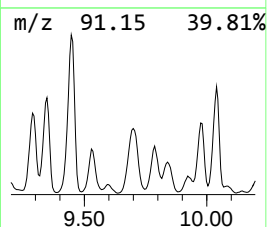
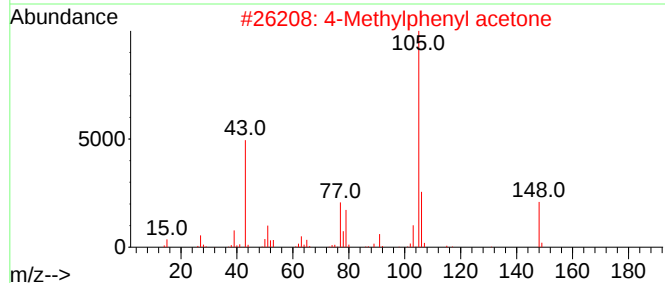
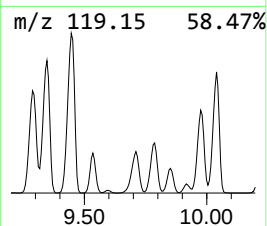
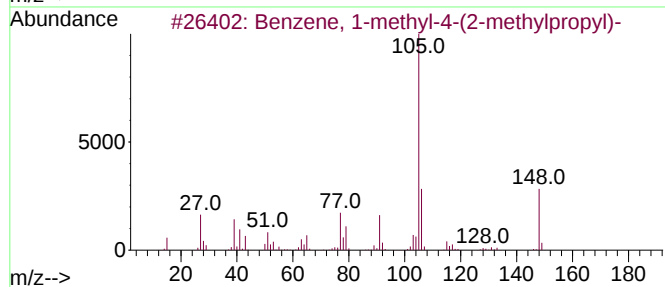
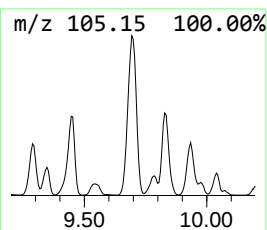
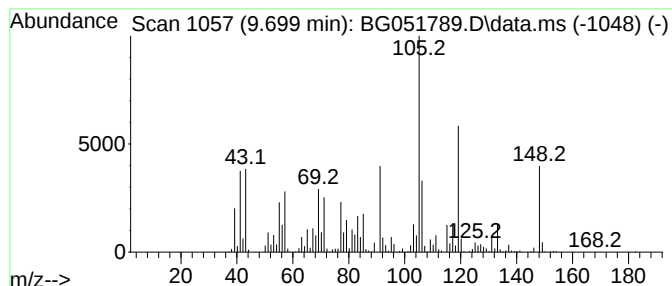
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 Benzene, 1-methyl-4-(2-meth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.699	12.57 ng/ul	13937600	1,4-Dichlorobenzene-d4	8.324

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	55
2			4-Methylphenyl acetone	148	C10H12O	002096-86-8	49
3			2-Methylindan-2-ol	148	C10H12O	033223-84-6	43
4			Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	41
5			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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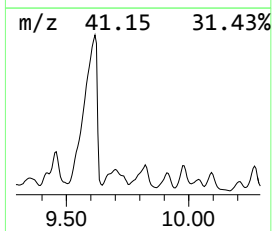
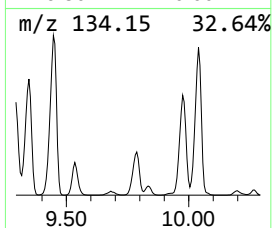
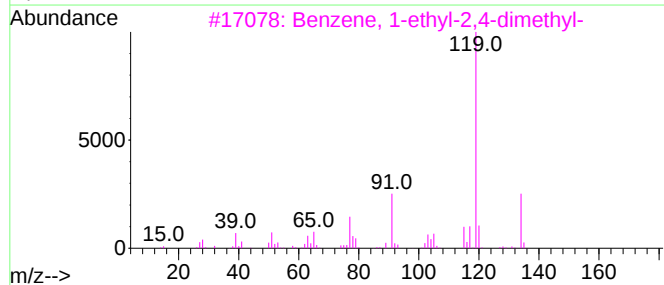
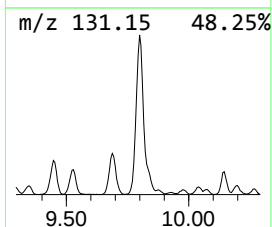
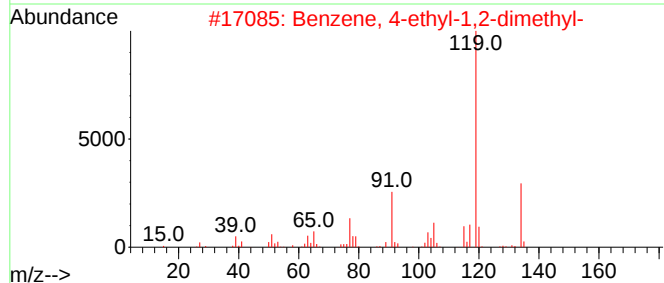
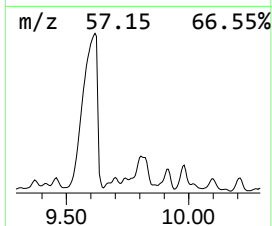
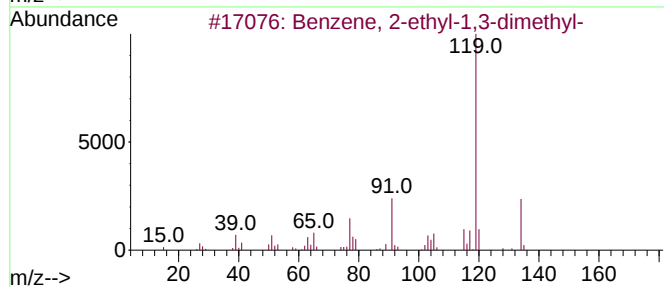
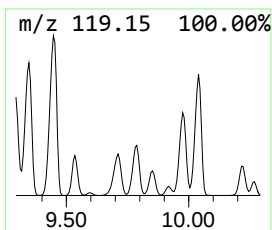
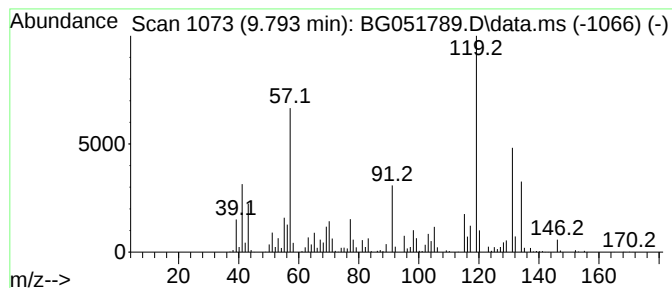
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.793	6.51 ng/ul	5409230	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	90
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	90
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	60
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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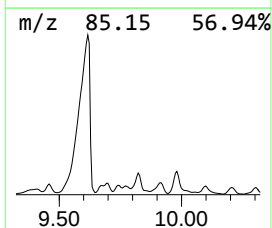
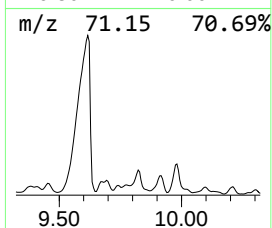
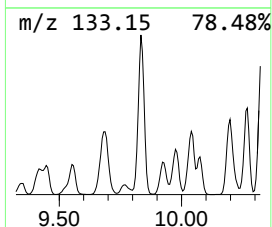
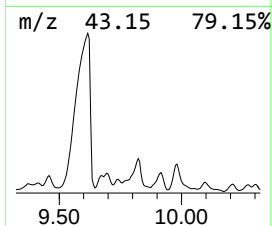
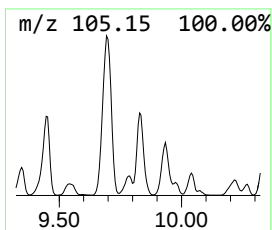
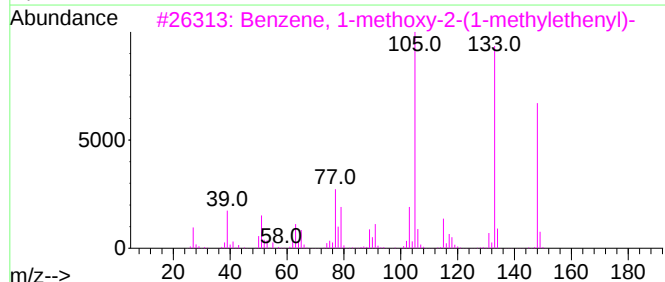
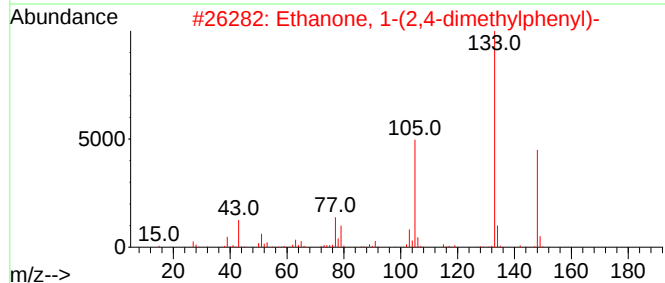
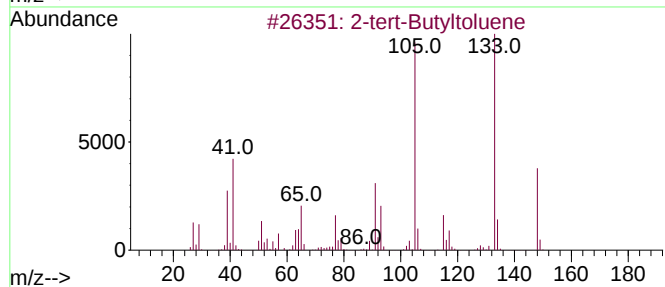
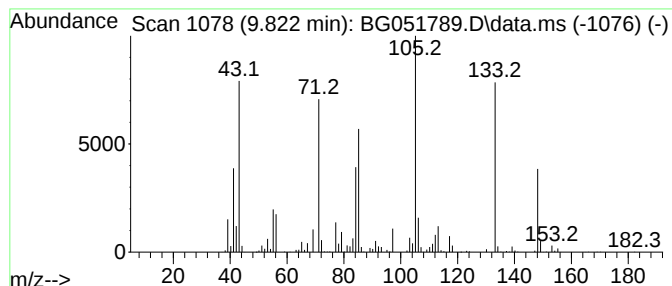
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 unknown-01 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.822	7.18 ng/ul	5960180	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-tert-Butyltoluene	148	C11H16	001074-92-6	43
2			Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	43
3			Benzene, 1-methoxy-2-(1-methylet...	148	C10H12O	010278-02-1	43
4			Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	41
5			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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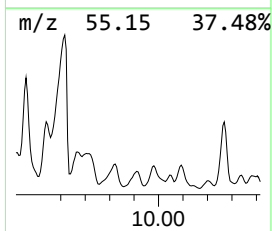
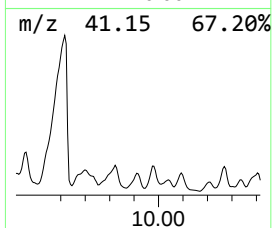
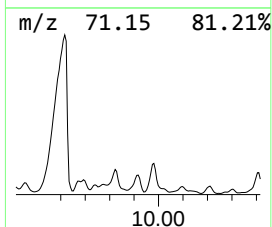
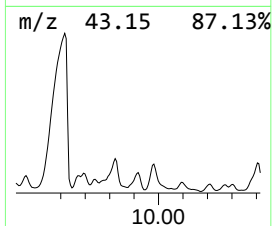
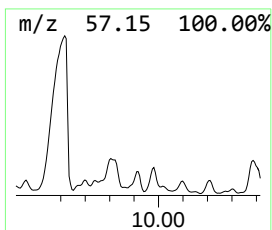
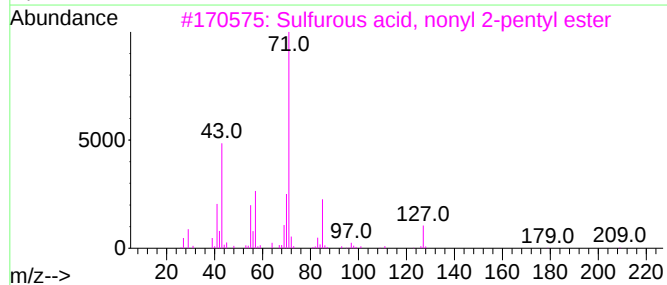
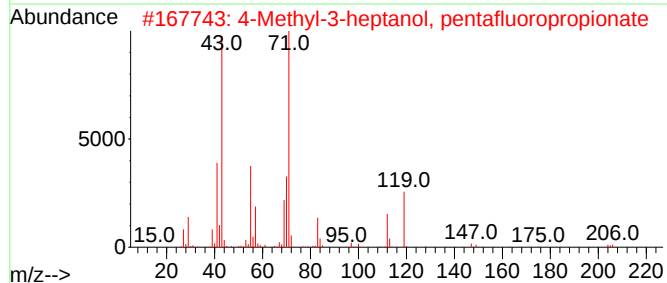
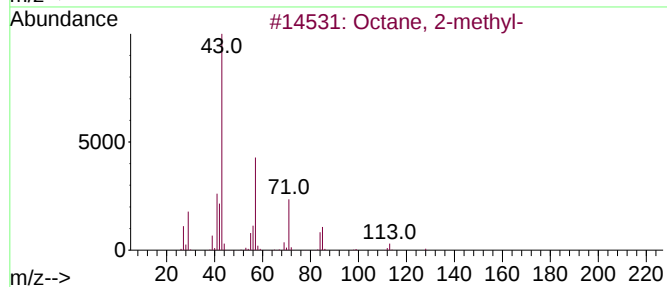
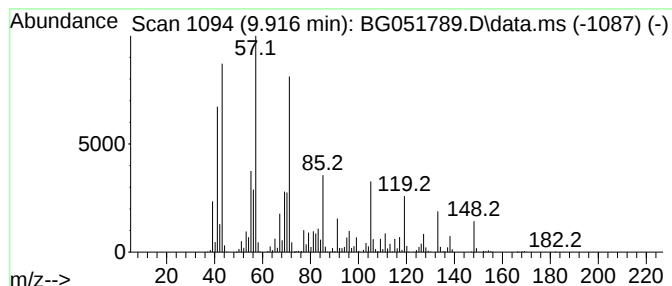
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.916	5.56 ng/ul	4612840	Naphthalene-d8	11.138

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2-methyl-	128	C9H20	003221-61-2	47
2		4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	46
3		Sulfurous acid, nonyl 2-pentyl e...	278	C14H30O3S	1010309-15-8	43
4		Octane, 3-ethyl-	142	C10H22	005881-17-4	43
5		Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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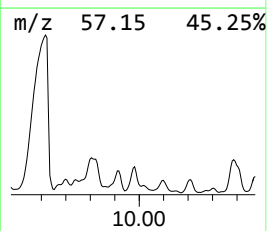
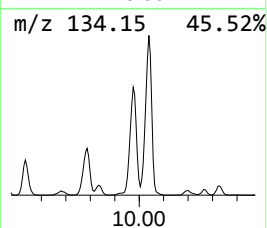
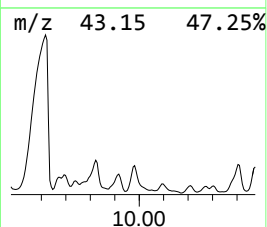
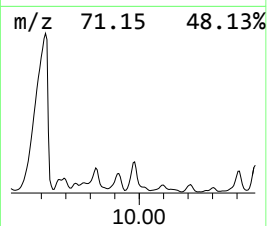
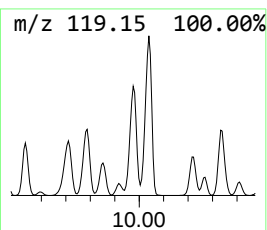
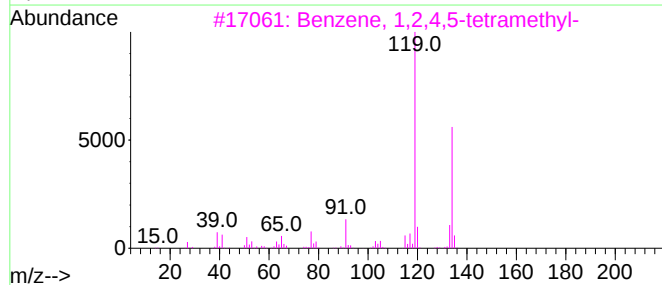
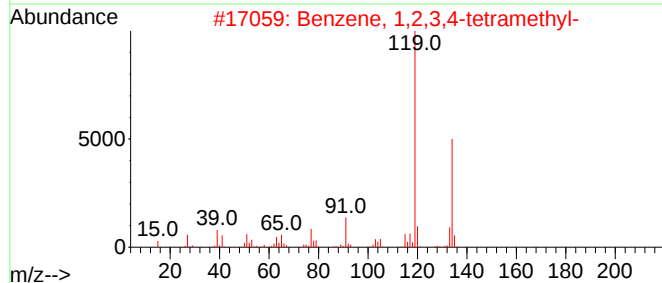
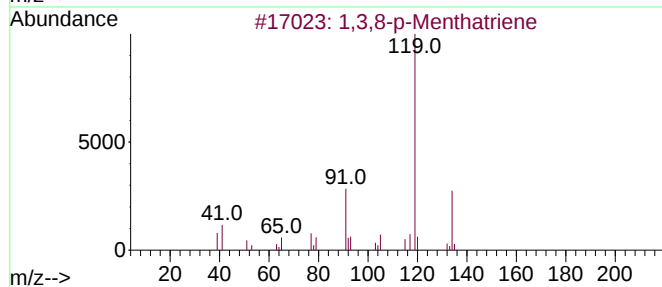
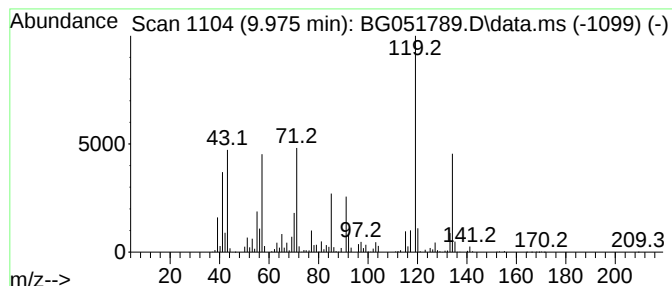
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 1,3,8-p-Menthatriene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.975	10.82 ng/ul	8981790	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	70
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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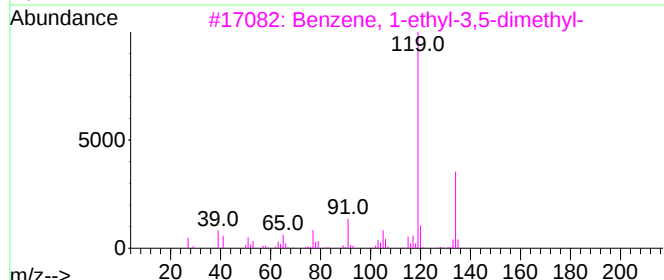
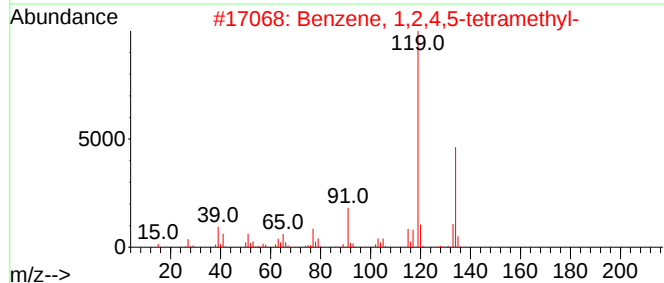
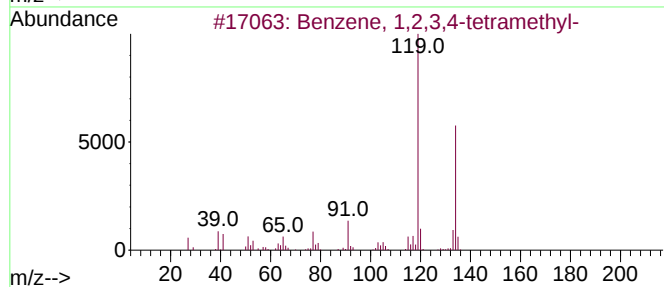
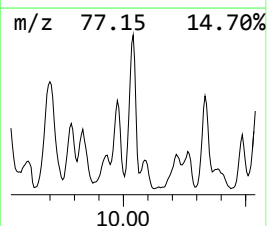
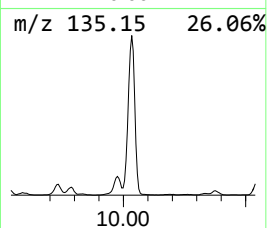
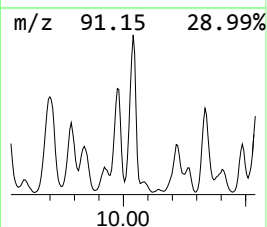
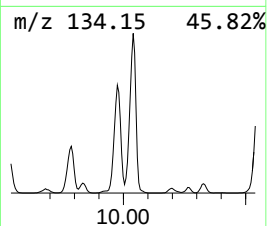
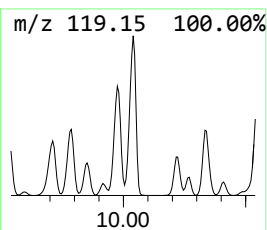
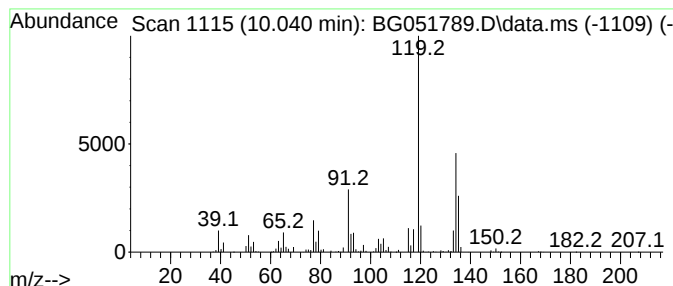
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.040	10.70 ng/ul	8884140	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	90
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
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Misc :
ALS Vial : 40 Sample Multiplier: 1

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ClientSampleId :
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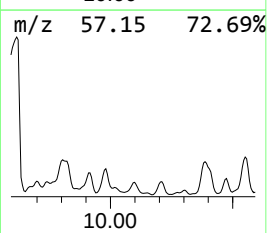
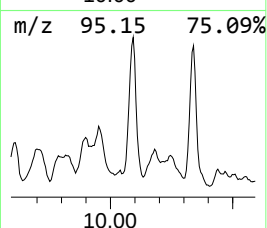
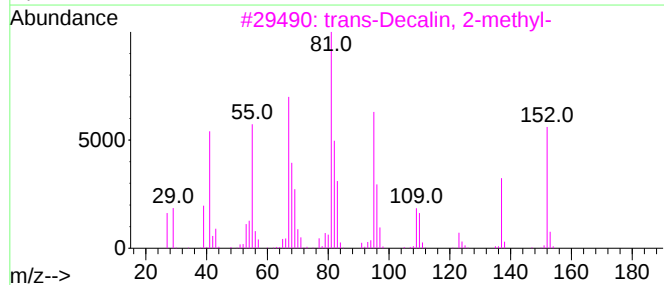
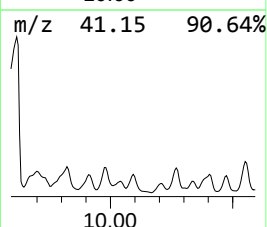
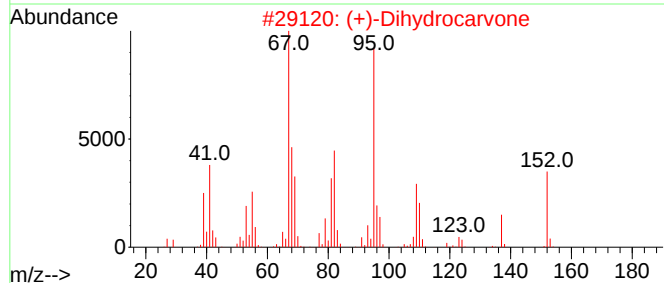
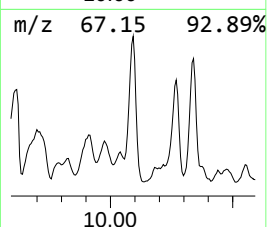
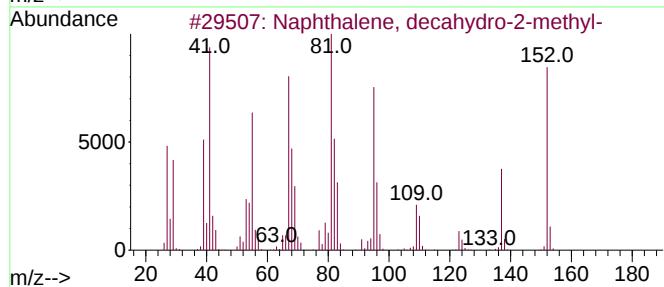
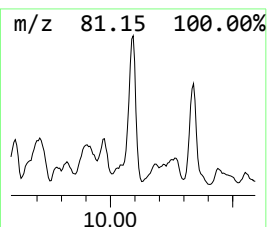
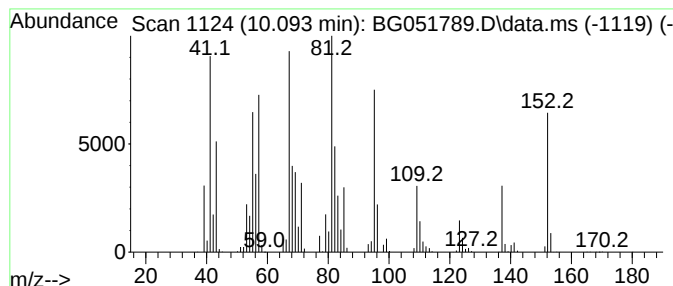
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 Naphthalene, decahydro-2-me... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.093	6.14 ng/ul	5100140	Naphthalene-d8	11.138

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	97
2		(+)-Dihydrocarvone	152	C10H16O	005524-05-0	89
3		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	87
4		Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	74
5		Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	64



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Data File : BG051789.D
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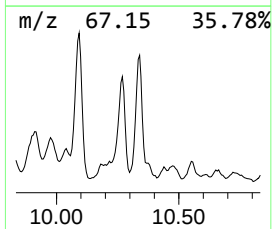
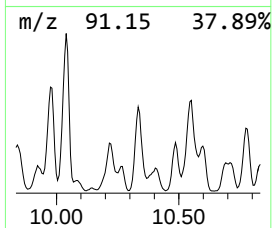
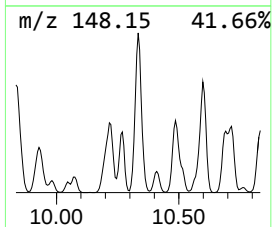
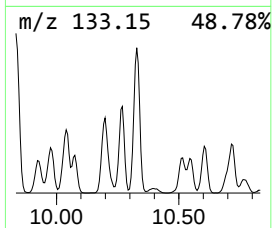
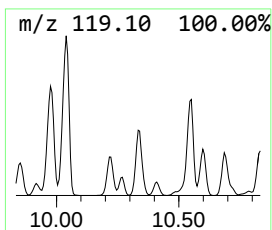
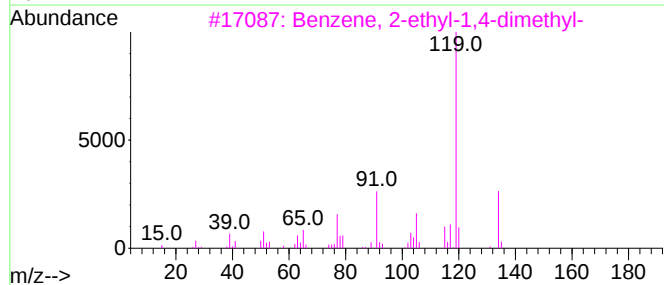
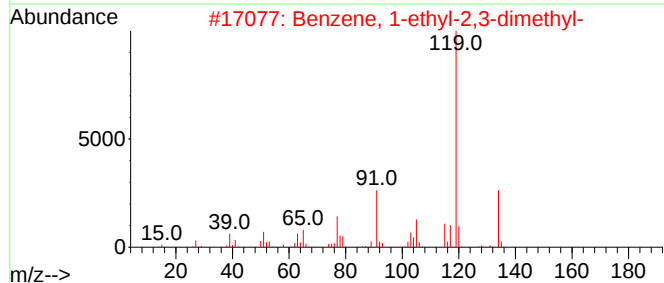
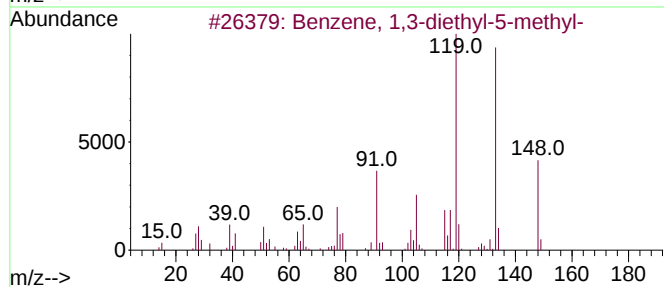
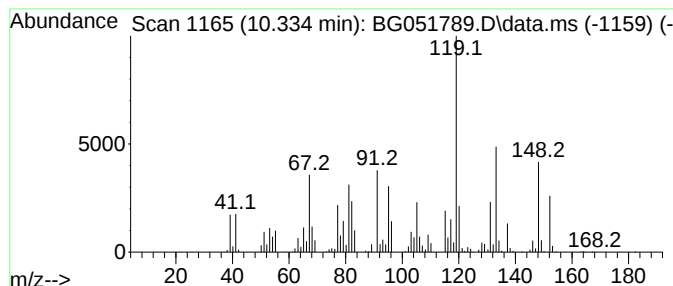
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.334	9.46 ng/ul	7856050	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	70
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	38
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
4			Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	38
5			3,4-Dihydro-2-[1H]quinoxalinone	148	C8H8N2O	059564-59-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

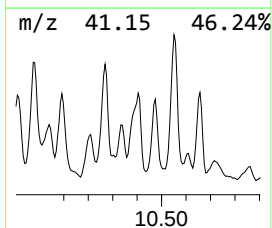
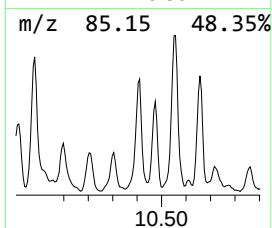
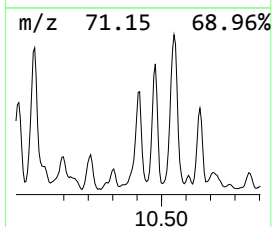
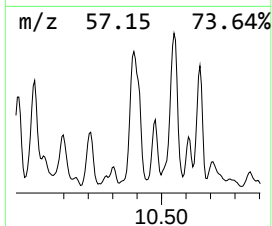
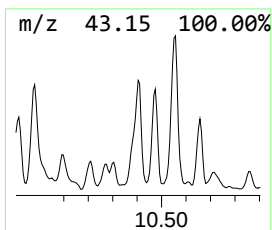
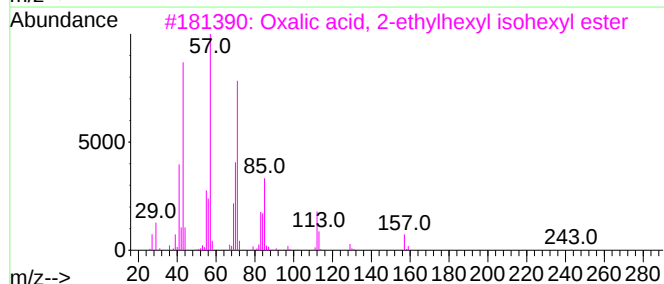
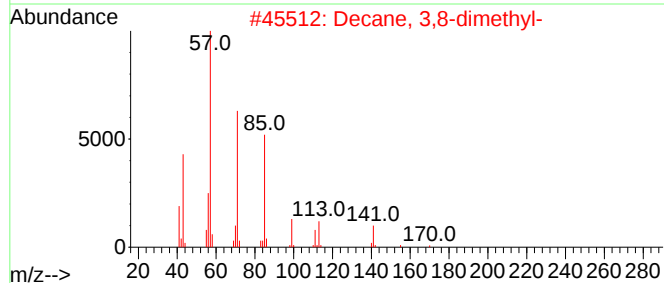
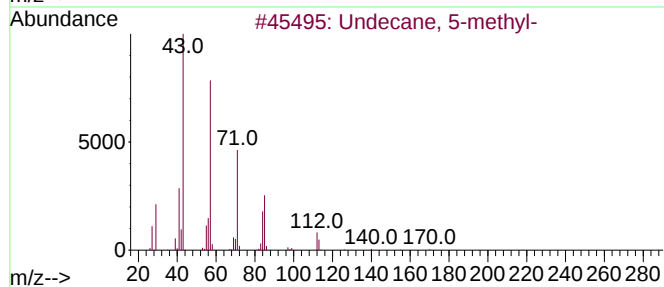
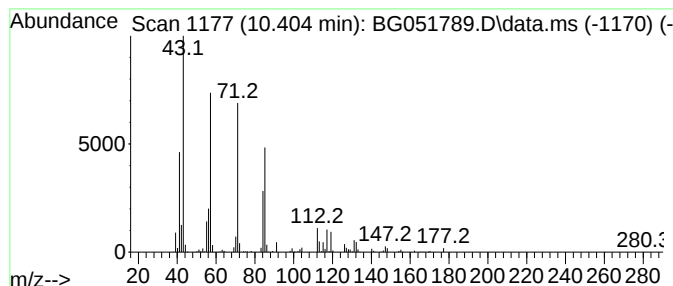
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.404	8.18 ng/ul	6793810	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 5-methyl-	170	C12H26	001632-70-8	58
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	52
3			Oxalic acid, 2-ethylhexyl isohex...	286	C16H30O4	1000309-38-8	47
4			Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	47
5			Nonane, 1-iodo-	254	C9H19I	004282-42-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

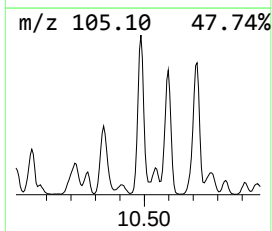
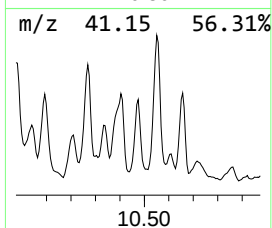
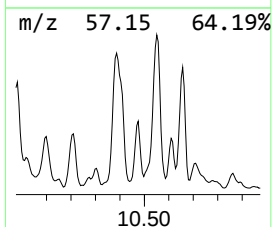
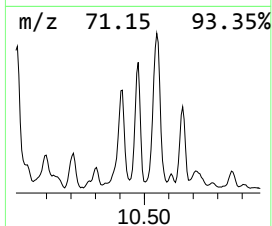
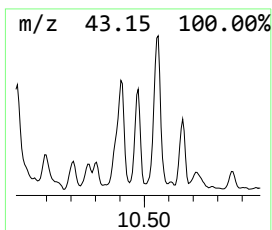
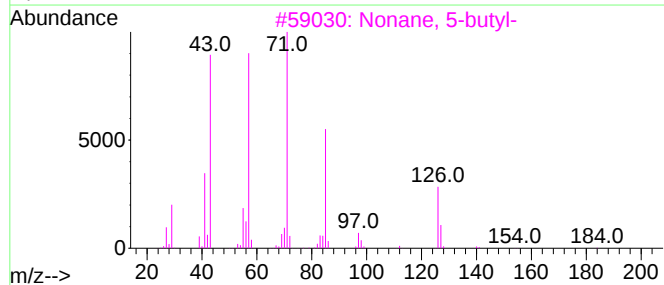
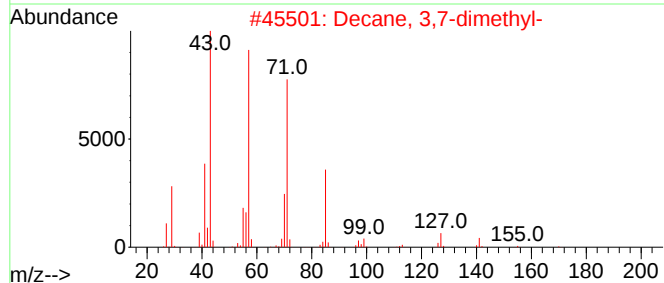
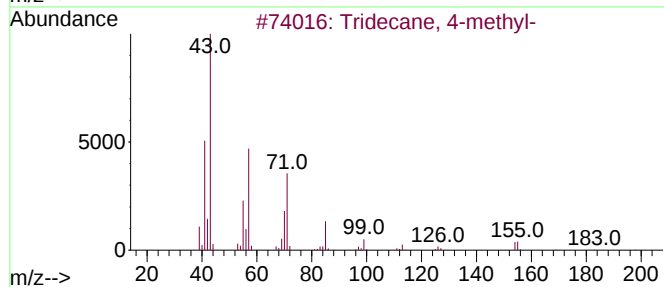
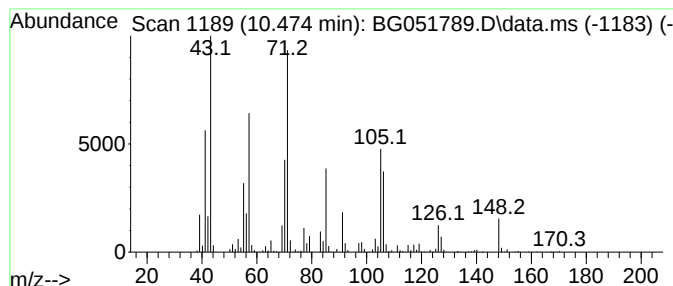
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.475	6.55 ng/ul	5437020	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 4-methyl-	198	C14H30	026730-12-1	53
2			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	50
3			Nonane, 5-butyl-	184	C13H28	017312-63-9	49
4			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	47
5			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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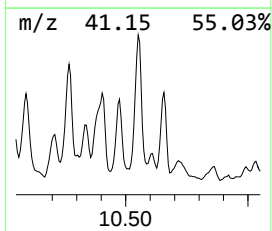
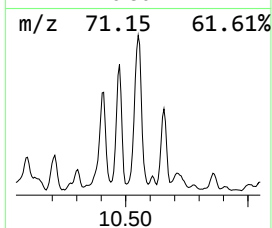
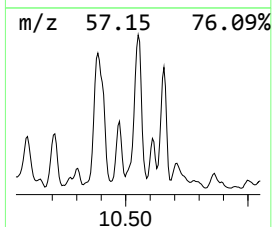
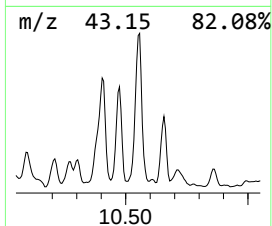
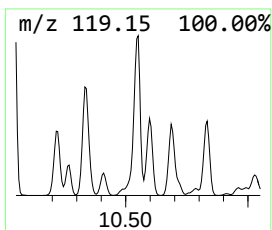
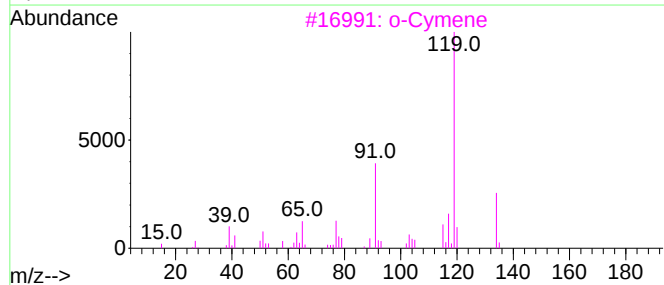
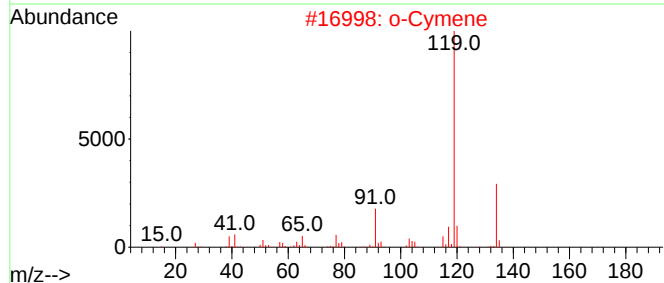
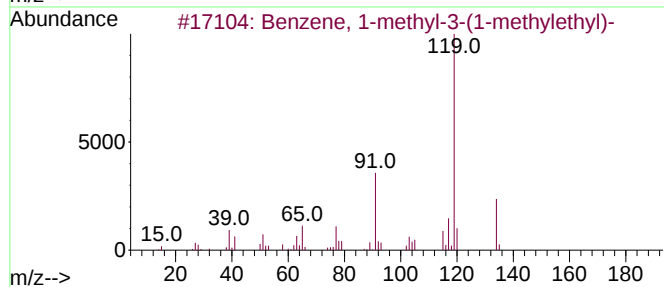
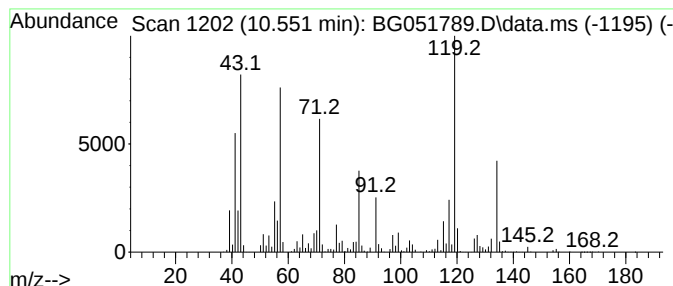
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Benzene, 1-methyl-3-(1-meth... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.551	14.26 ng/ul	11838100	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	55
2			o-Cymene	134	C10H14	000527-84-4	55
3			o-Cymene	134	C10H14	000527-84-4	55
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	50
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

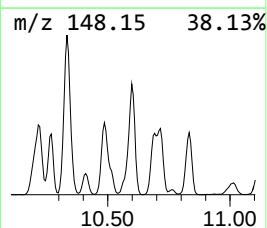
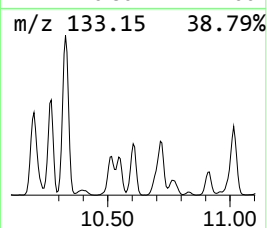
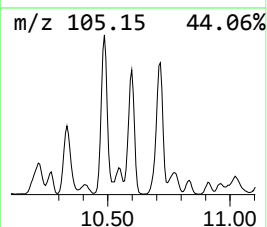
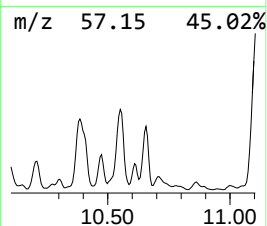
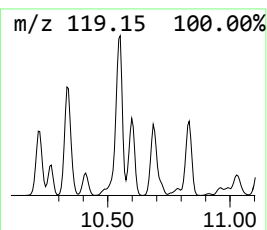
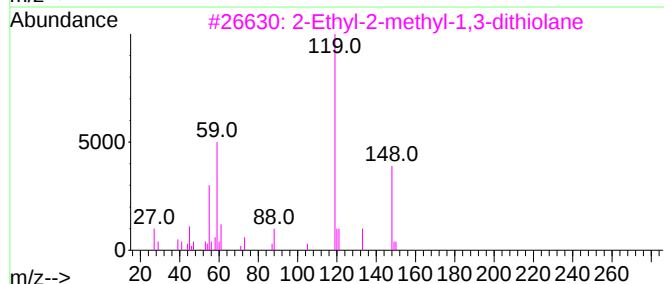
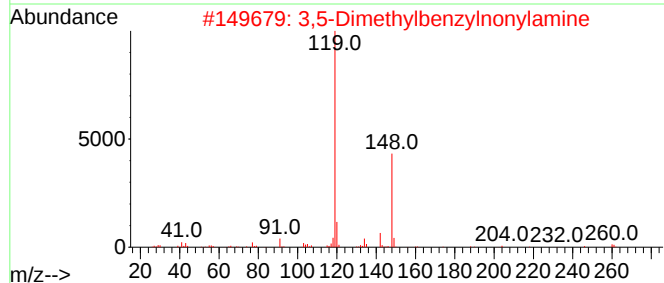
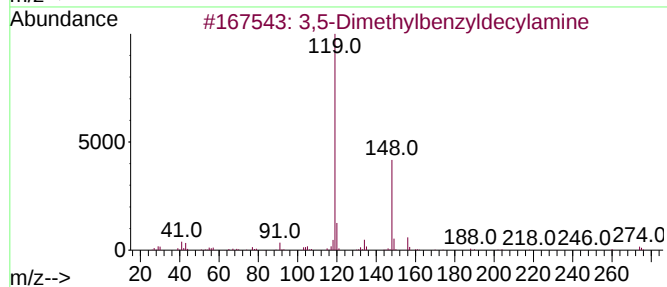
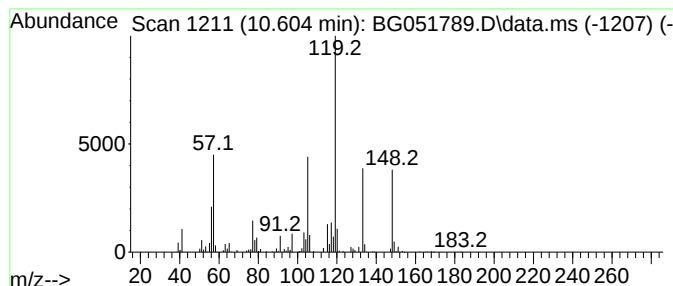
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 3,5-Dimethylbenzyldecylamine Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	3.55 ng/ul	2947370	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dimethylbenzyldecylamine	275	C19H33N	1000491-61-2	50
2			3,5-Dimethylbenzylnonylamine	261	C18H31N	1010491-61-1	50
3			2-Ethyl-2-methyl-1,3-dithiolane	148	C6H12S2	006008-81-7	49
4			3,4-Dihydro-2-[1H]quinoxalinone	148	C8H8N2O	059564-59-9	47
5			Benzoic acid, 2,4-dimethyl-, (2,...	268	C18H20O2	055000-43-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

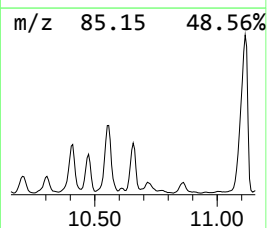
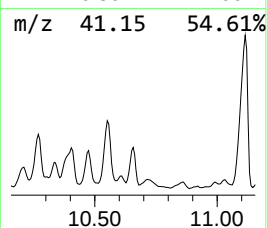
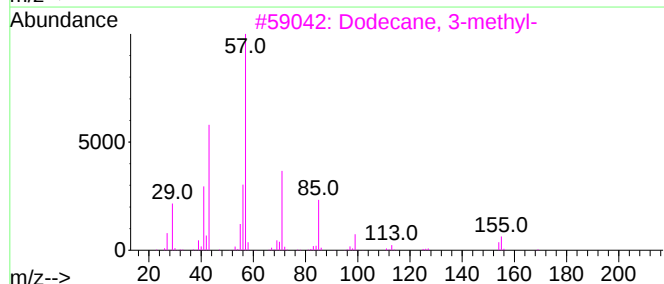
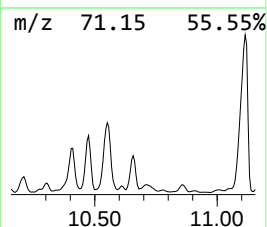
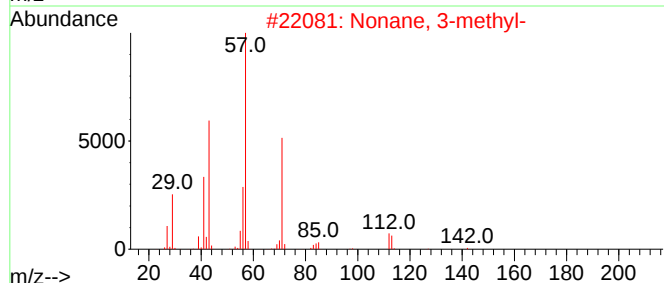
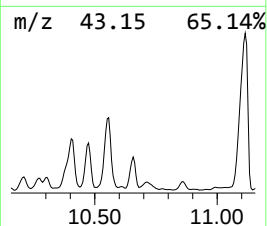
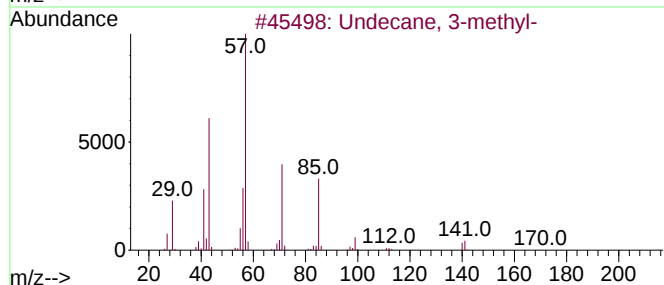
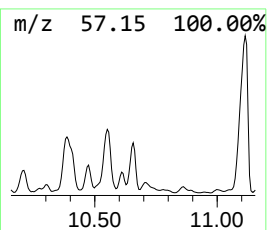
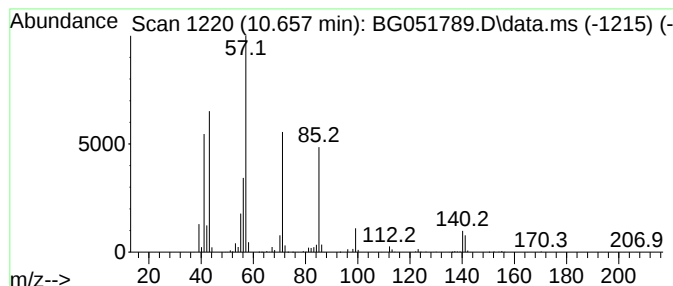
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.657	3.75 ng/ul	3110350	Naphthalene-d8	11.138

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3-methyl-	170	C12H26	001002-43-3	86
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	53
3		Dodecane, 3-methyl-	184	C13H28	017312-57-1	53
4		Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	53
5		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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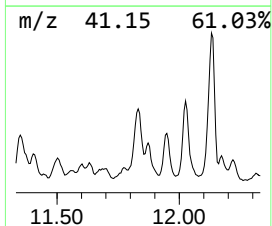
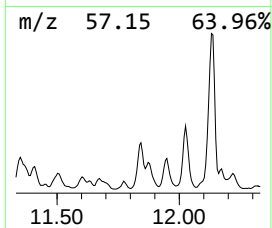
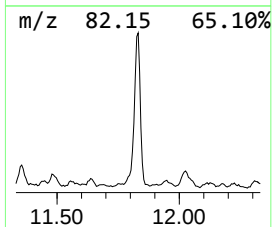
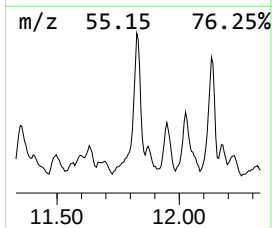
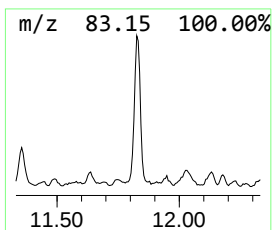
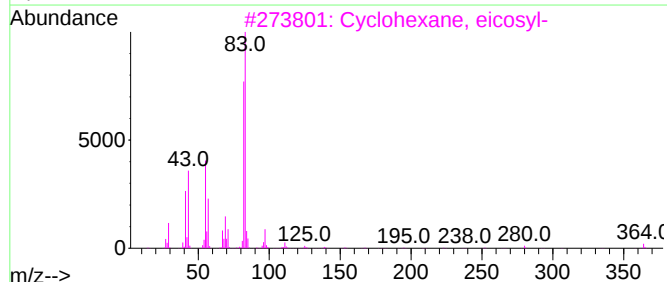
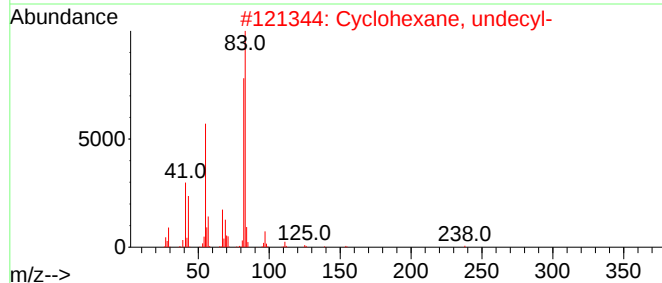
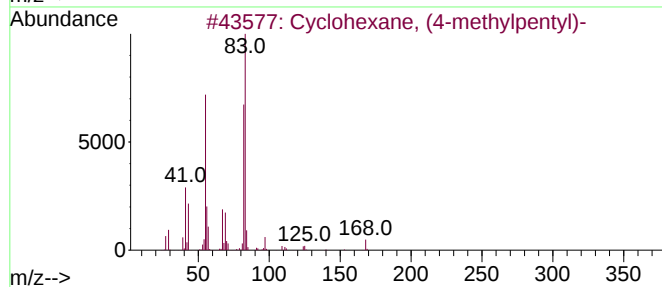
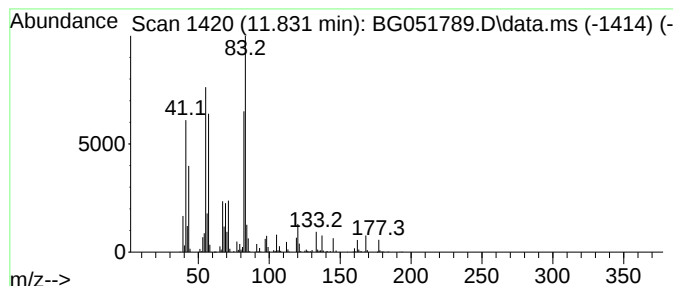
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Cyclic11.832 Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.832	2.24 ng/ul	1857150	Naphthalene-d8	11.138

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	81
2		Cyclohexane, undecyl-	238	C17H34	054105-66-7	64
3		Cyclohexane, eicosyl-	364	C26H52	004443-55-4	59
4		Cyclohexane, pentyl-	154	C11H22	004292-92-6	59
5		Cyclohexane, octyl-	196	C14H28	001795-15-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

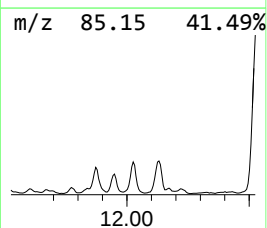
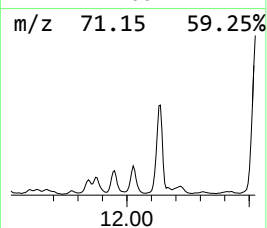
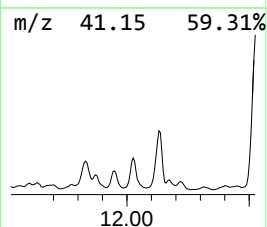
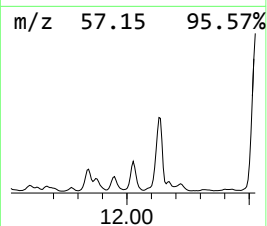
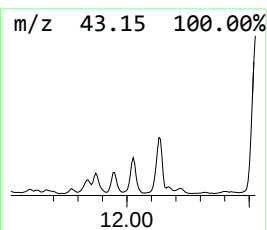
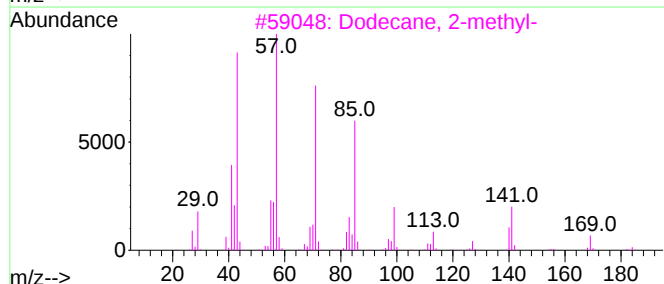
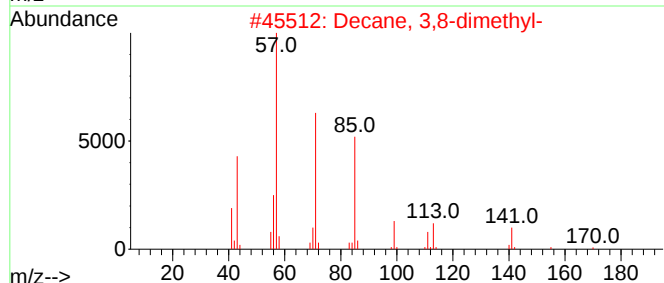
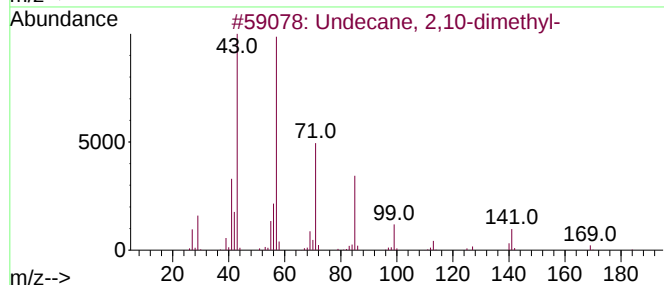
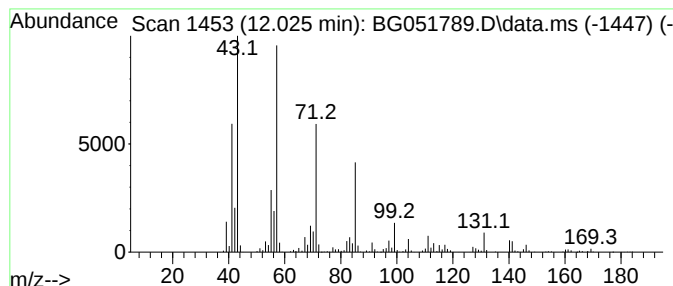
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.025	2.50 ng/ul	2074380	Naphthalene-d8	11.138

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 2,10-dimethyl-	184	C13H28	017301-27-8	91
2		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	76
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	70
4		Decane, 1-iodo-	268	C10H21I	002050-77-3	64
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

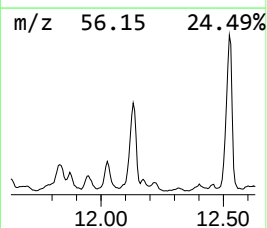
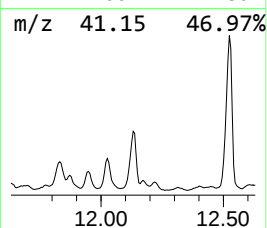
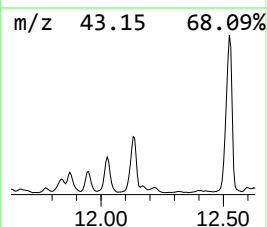
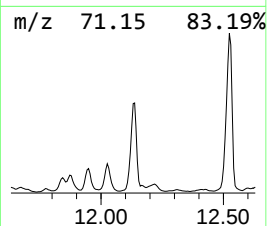
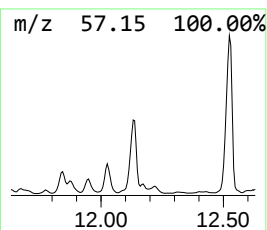
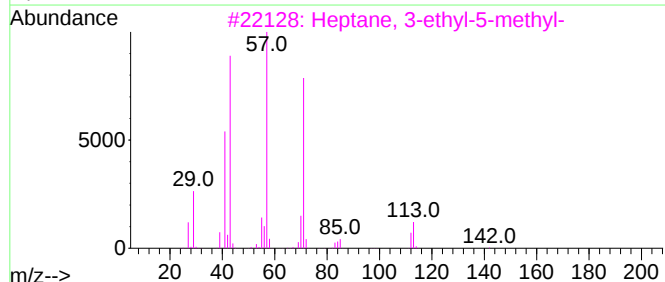
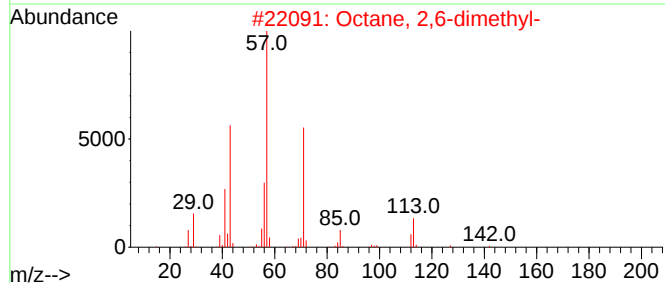
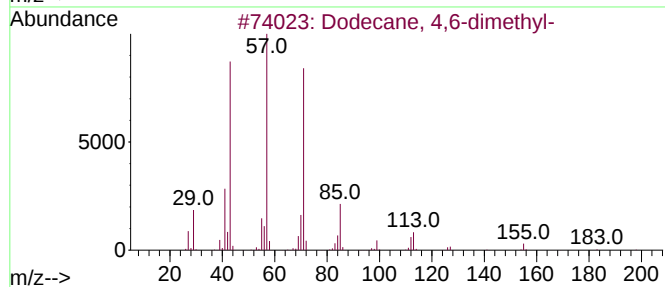
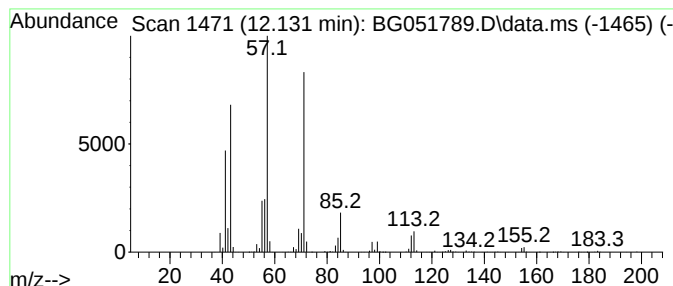
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.131	4.60 ng/ul	3819000	Naphthalene-d8	11.138

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	91
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
3			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	80
4			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

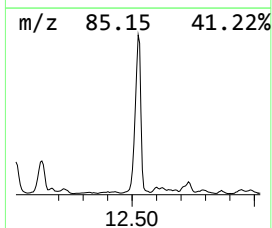
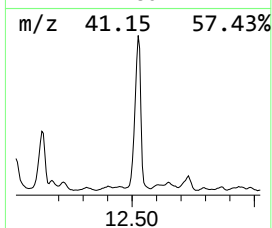
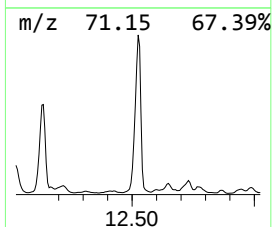
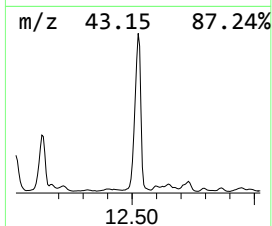
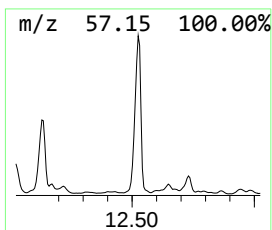
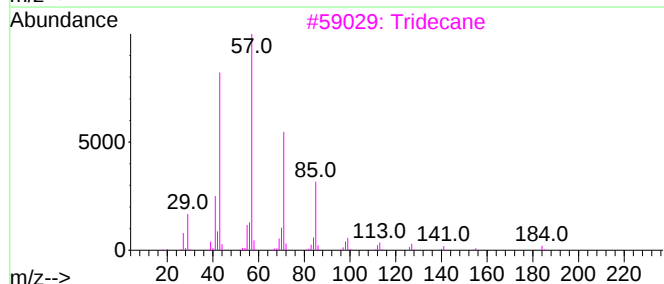
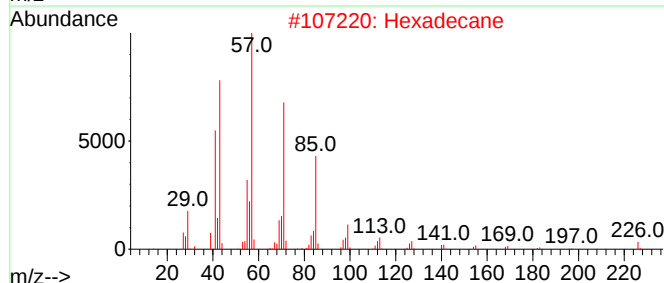
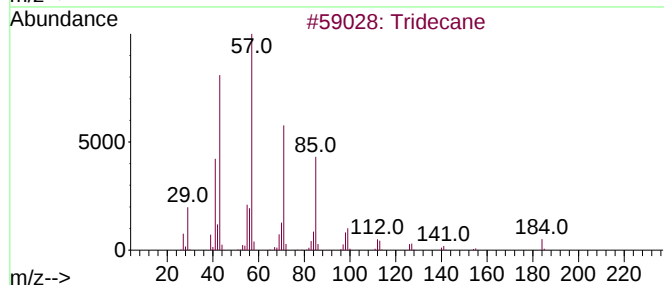
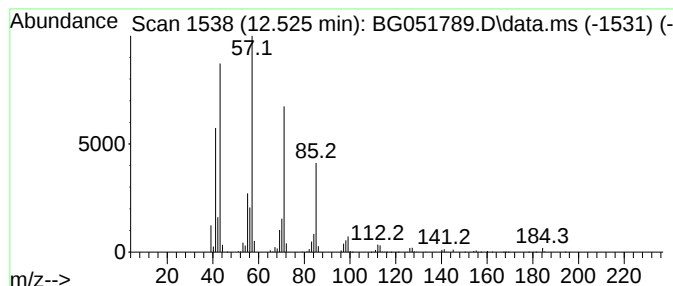
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.525	10.23 ng/ul	8492810	Naphthalene-d8	11.138	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	97
2	Hexadecane	226	C16H34	000544-76-3	91
3	Tridecane	184	C13H28	000629-50-5	90
4	Tetradecane	198	C14H30	000629-59-4	90
5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

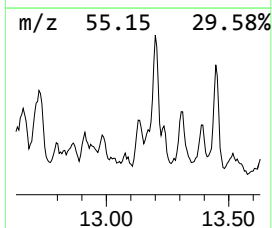
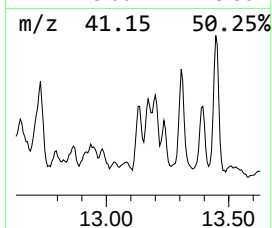
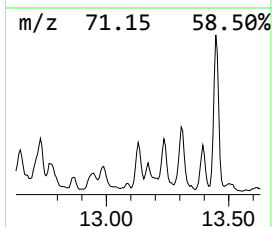
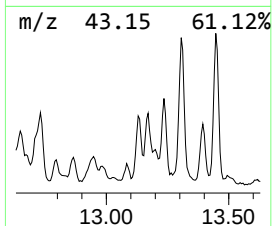
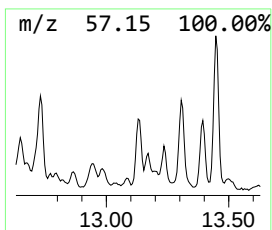
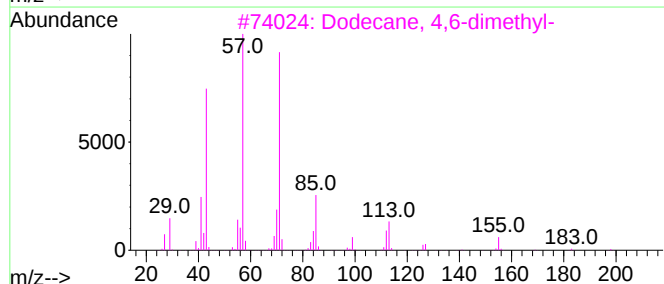
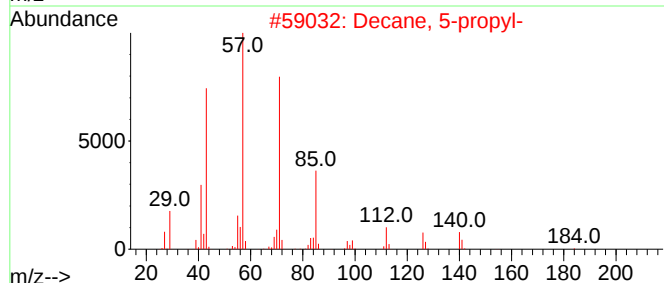
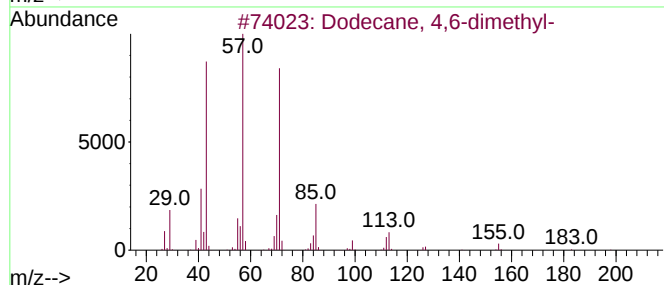
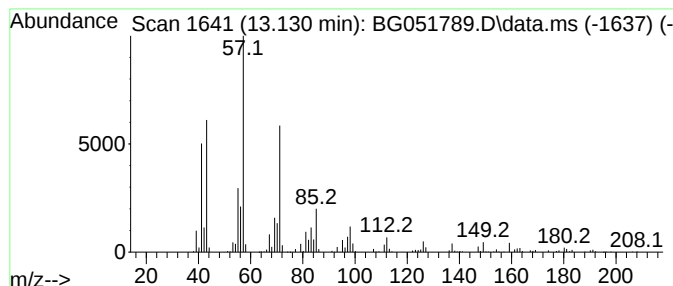
Instrument :
BNA_G
ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.130	13.33 ng/ul	690663	Acenaphthene-d10	14.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64
2	Decane, 5-propyl-	184	C13H28	017312-62-8	50
3	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	47
4	Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	47
5	2,3-Dimethyldodecane	198	C14H30	006117-98-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

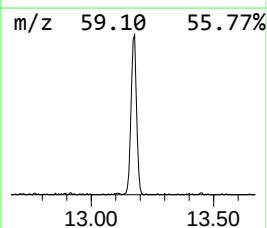
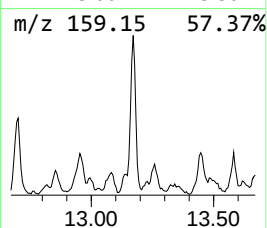
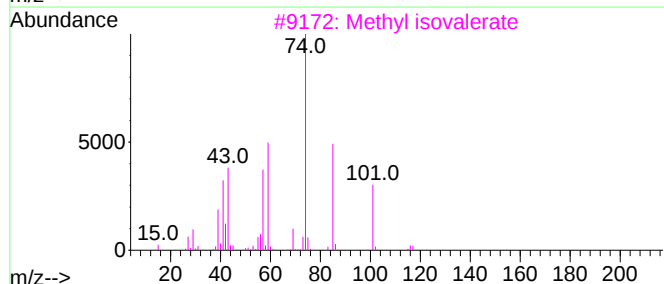
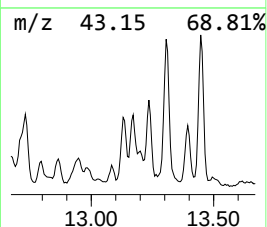
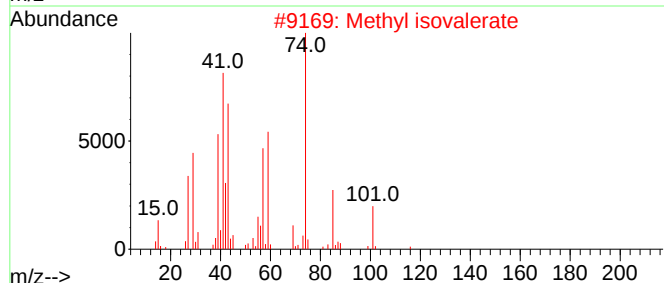
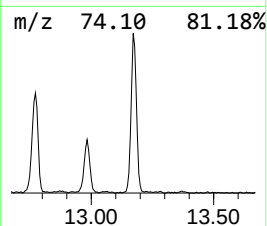
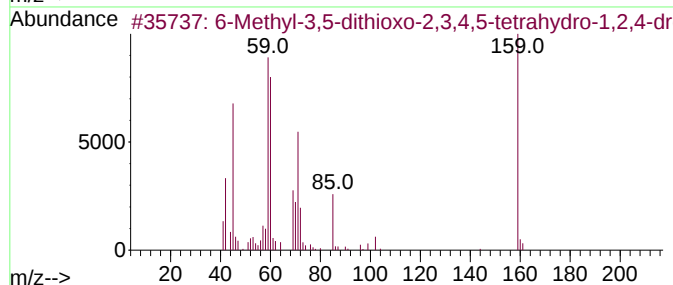
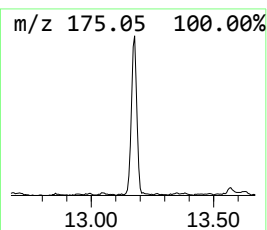
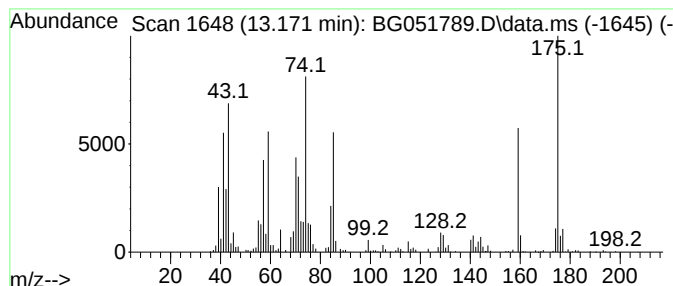
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 unknown-02 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.171	11.95 ng/ul	619273	Acenaphthene-d10		14.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-3,5-dithioxo-2,3,4,5-te...	159	C4H5N3S2	000452-88-0	14	
2	Methyl isovalerate	116	C6H12O2	000556-24-1	14	
3	Methyl isovalerate	116	C6H12O2	000556-24-1	12	
4	Allyldimethyl(vinyl)silane	126	C7H14Si	022146-25-4	10	
5	4-quinolinethiol, 2-methyl-	175	C10H9NS	1000404-71-3	10	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

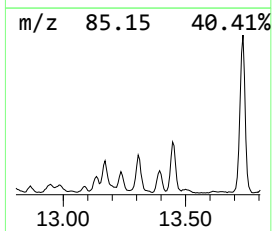
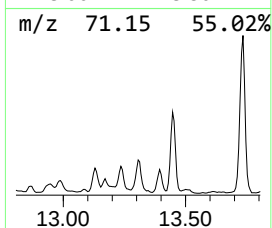
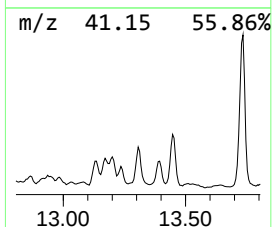
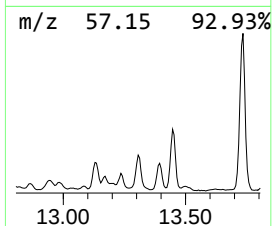
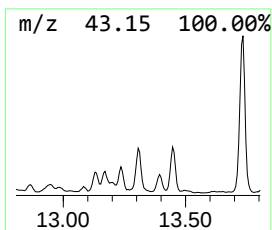
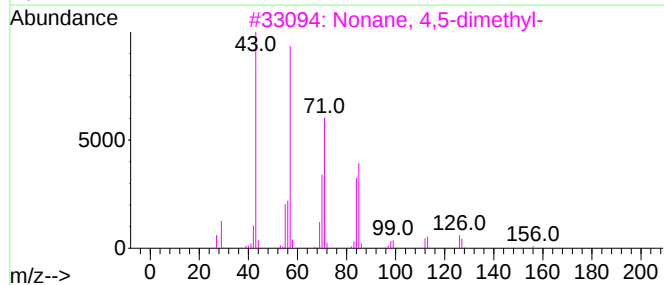
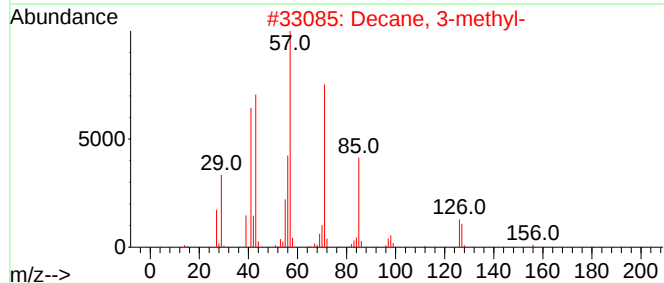
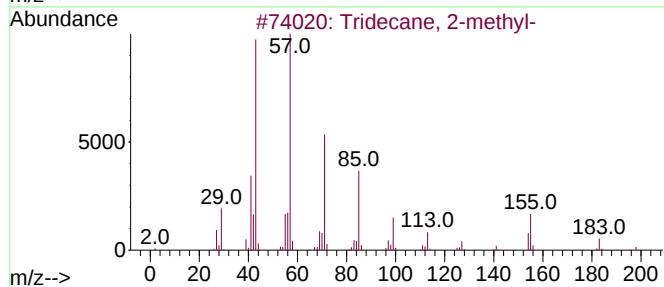
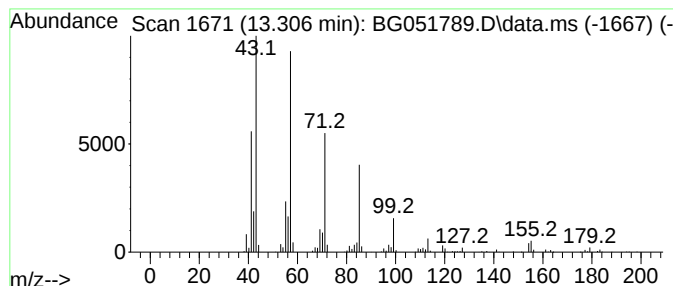
Instrument :
BNA_G
ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.306	16.09 ng/ul	833779	Acenaphthene-d10	14.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 2-methyl-	198	C14H30	001560-96-9	80
2	Decane, 3-methyl-	156	C11H24	013151-34-3	72
3	Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	72
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	72
5	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

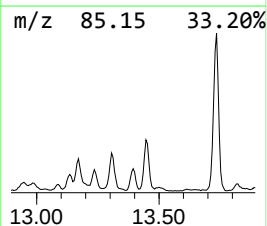
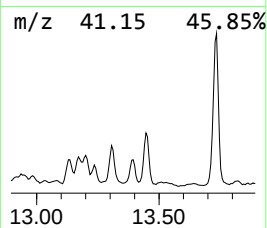
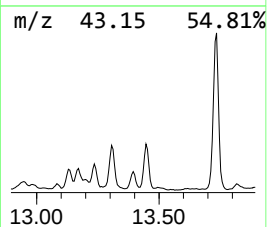
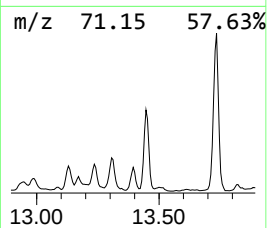
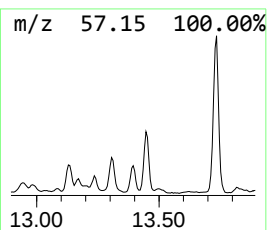
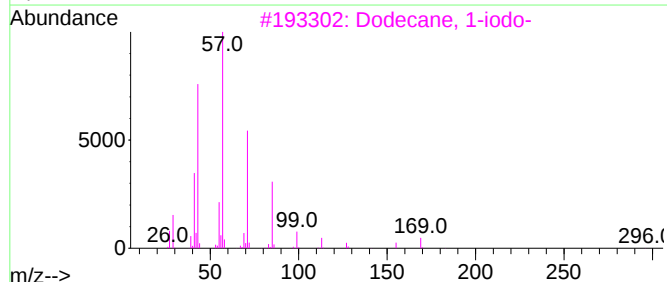
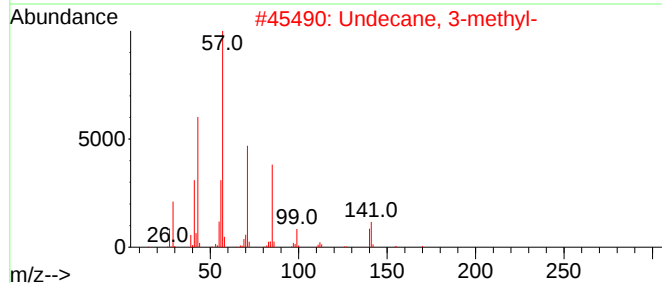
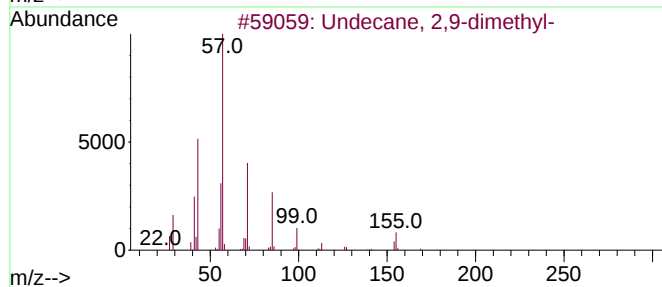
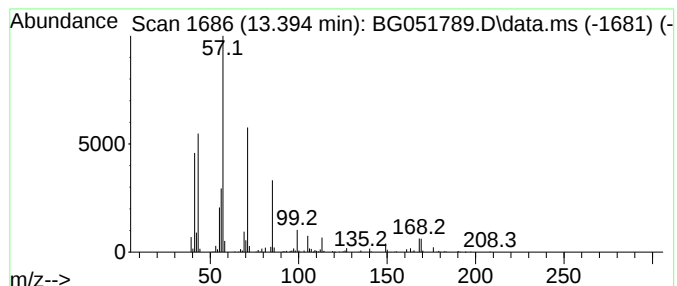
Instrument :
BNA_G
ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.394	10.24 ng/ul	530638	Acenaphthene-d10		14.916	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 2,9-dimethyl-		184	C13H28	017301-26-7	78
2	Undecane, 3-methyl-		170	C12H26	001002-43-3	72
3	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	72
4	Tridecane, 3-methyl-		198	C14H30	006418-41-3	58
5	Undecane, 5-methyl-		170	C12H26	001632-70-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

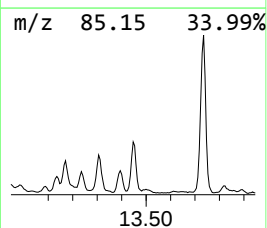
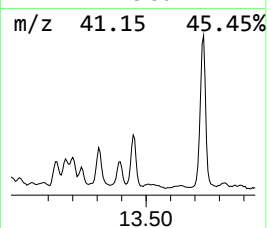
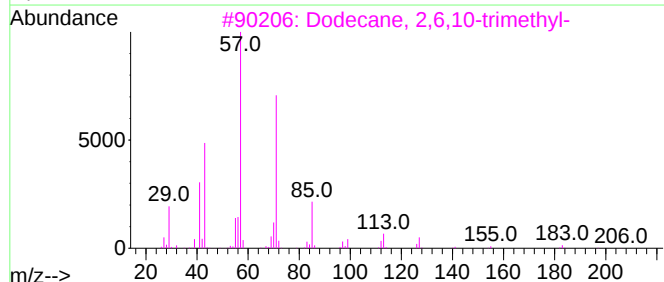
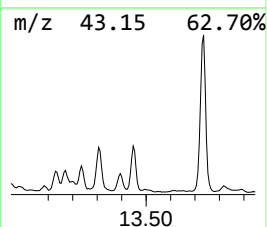
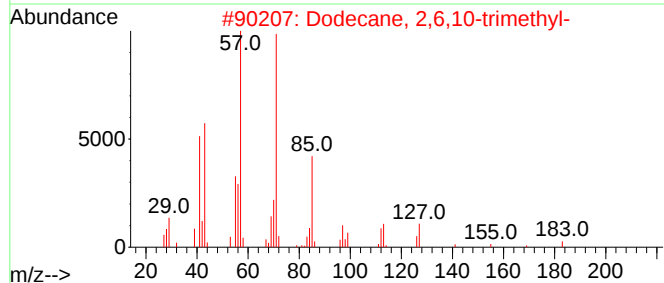
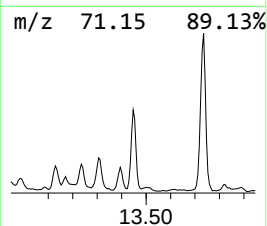
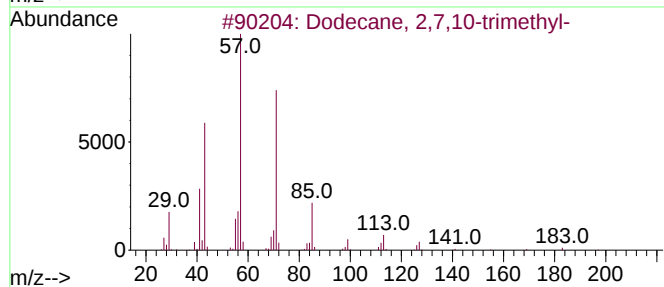
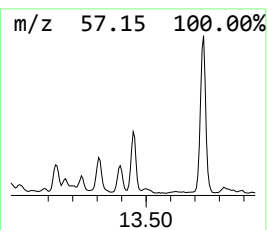
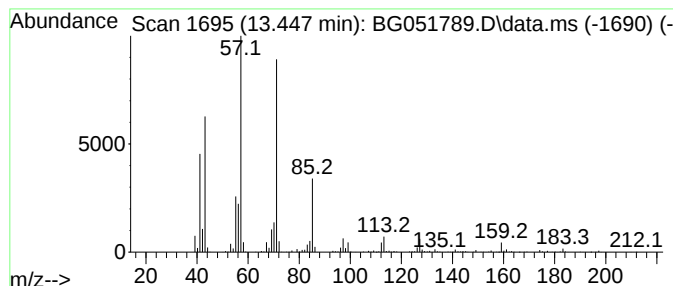
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Quant Title : SVOA CALIBRATION

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

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	26.49 ng/ul	1373000	Acenaphthene-d10	14.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	90
2			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	90
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	86
4			Decane, 5-propyl-	184	C13H28	017312-62-8	74
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

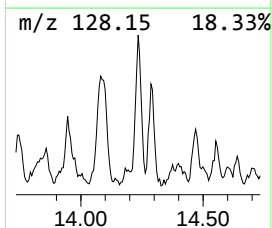
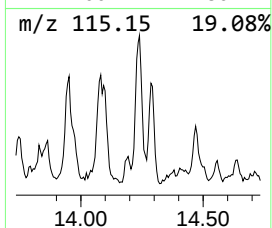
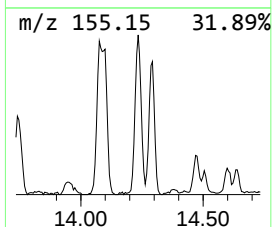
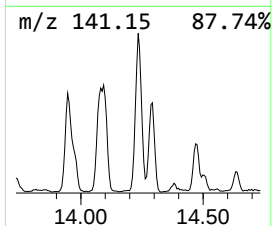
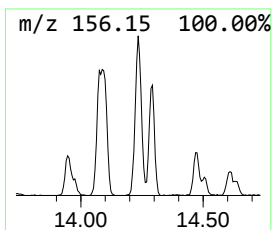
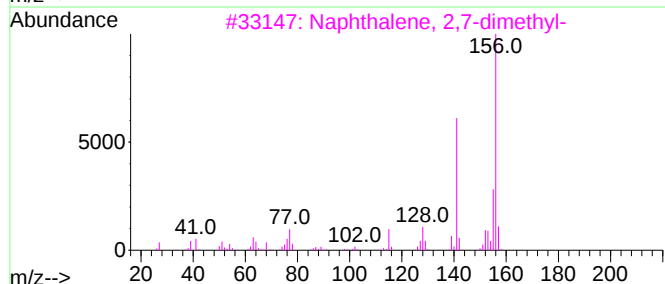
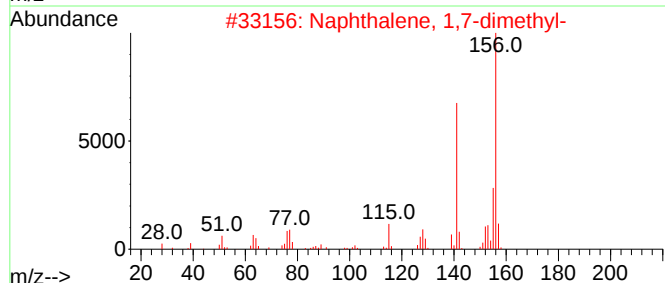
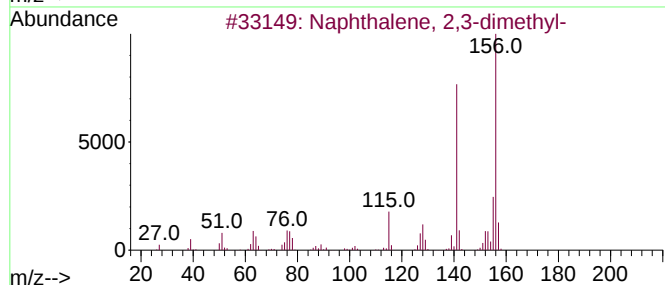
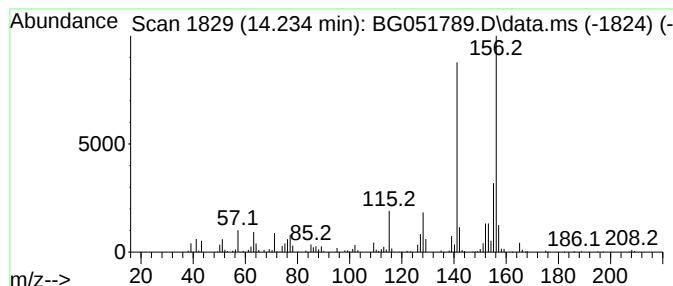
Instrument :
BNA_G
ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 56 Naphthalene, 2,3-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.234	11.83 ng/ul	613330	Acenaphthene-d10	14.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

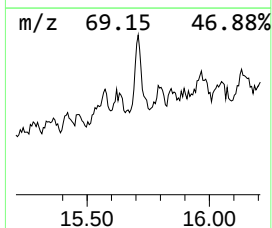
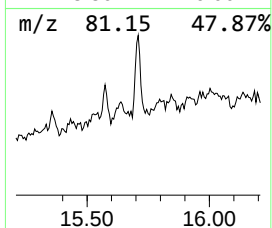
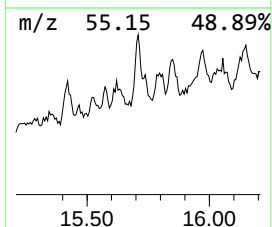
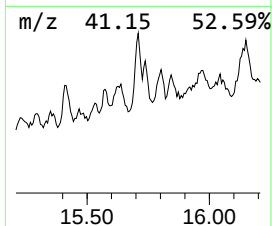
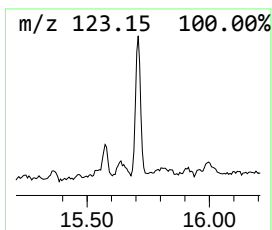
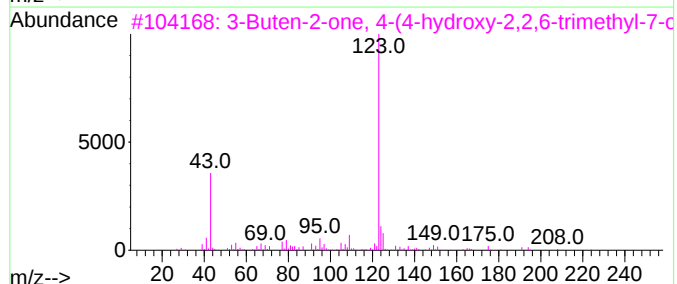
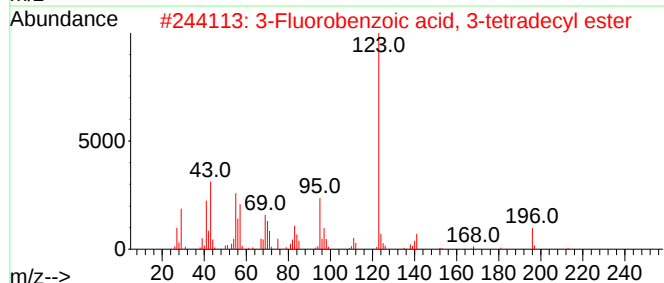
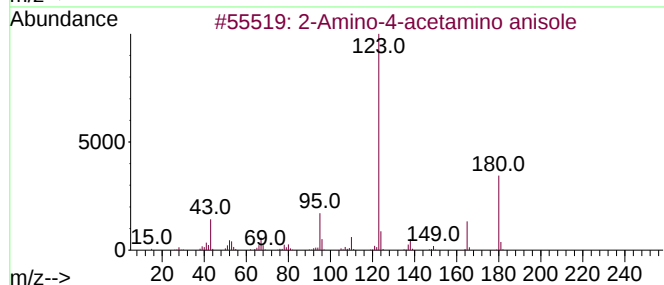
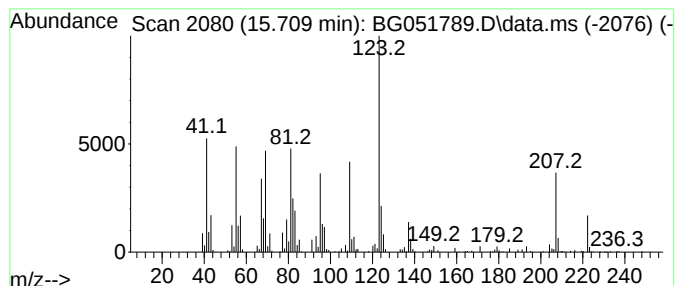
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 unknown-03 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.709	10.19 ng/ul	527911	Acenaphthene-d10	14.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Amino-4-acetamino anisole	180	C9H12N2O2	006375-47-9	38
2		3-Fluorobenzoic acid, 3-tetradec...	336	C21H33FO2	1000280-60-5	38
3		3-Buten-2-one, 4-(4-hydroxy-2,2,...	224	C13H20O3	038274-01-0	38
4		3-(2-Ethyl-imidazol-1-yl)-propio...	168	C8H12N2O2	1010277-78-0	38
5		4(1H)-Pyrimidinone, 6-(1-methyle...	180	C9H12N2O2	1000503-76-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

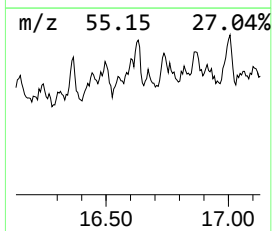
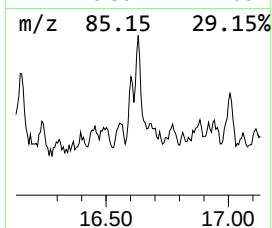
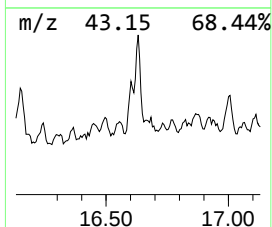
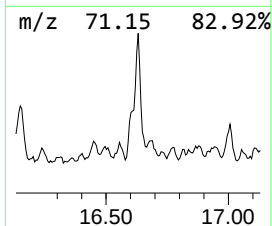
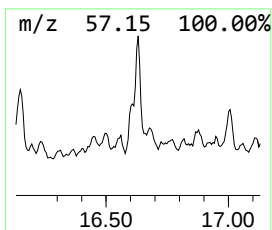
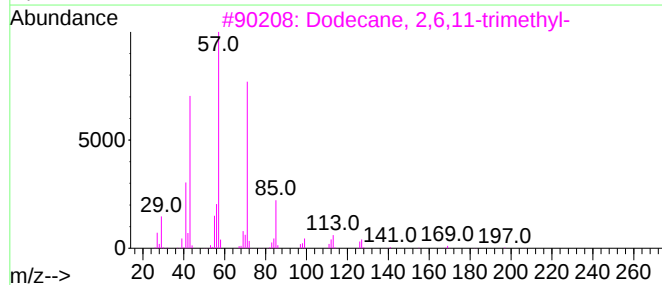
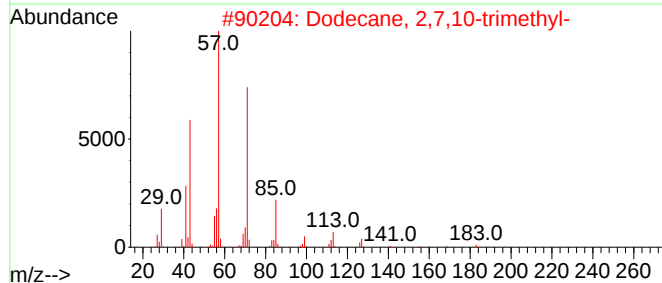
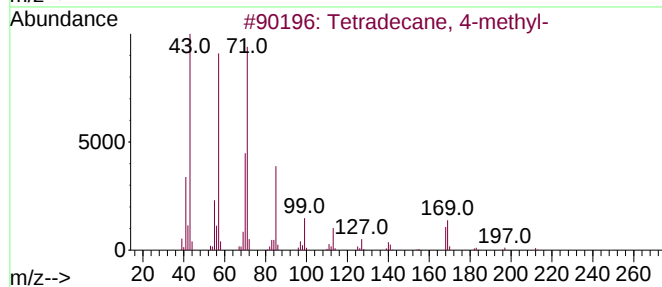
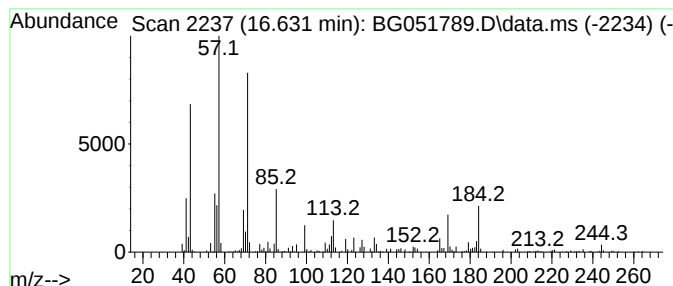
Instrument :
BNA_G
ClientSampleId :
BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
16.631	20.00 ng/ul	623194	Phenanthrene-d10			17.665
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane, 4-methyl-		212	C15H32	025117-24-2	76
2	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	59
3	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59
4	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	58
5	Heptacosane		380	C27H56	000593-49-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
Data File : BG051789.D
Acq On : 23 Dec 2021 17:44
Operator : CG/JU
Sample : M5215-04
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD6

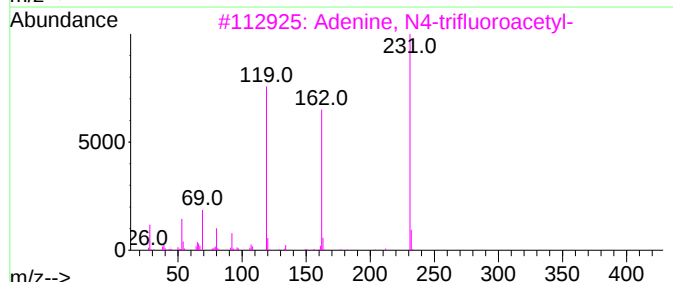
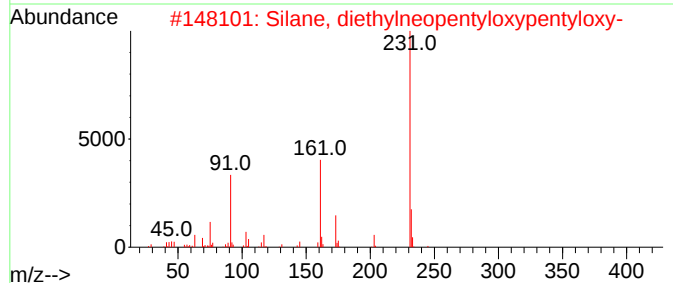
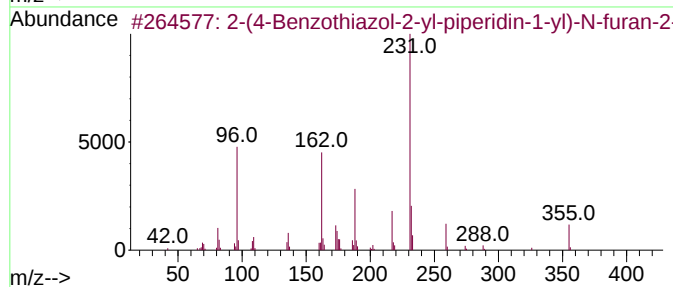
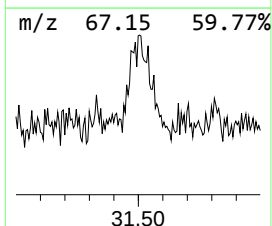
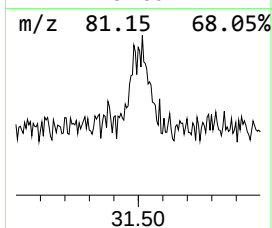
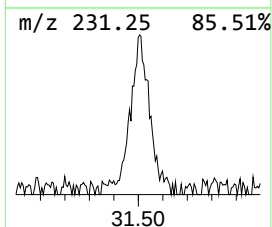
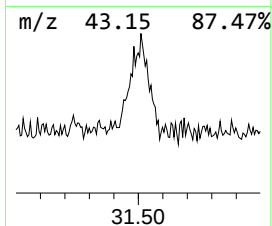
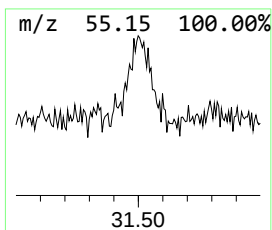
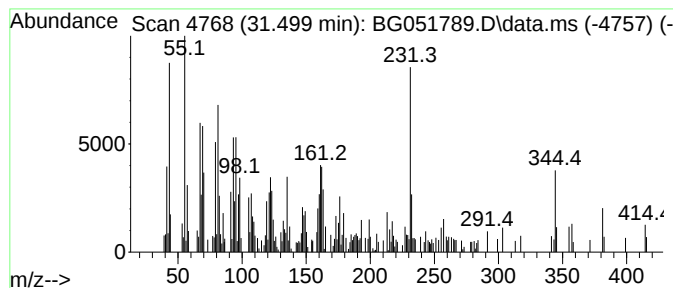
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 59 unknown-04 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.499	19.72 ng/ul	502085	Perylene-d12	25.466

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-Benzothiazol-2-yl-piperidin...	355	C19H21N3O2S	1000322-21-0	41		
2	Silane, diethyneopentyloxy...	260	C14H32O2Si	1000363-94-0	35		
3	Adenine, N4-trifluoroacetyl-	231	C7H4F3N5O	1000374-88-1	35		
4	Benzoic acid, 2-(5-pyrazolylcarb...	231	C11H9N3O3	1000277-53-0	25		
5	Pyrimidine, 4-amino-2,6-bis(trif...	231	C6H3F6N3	000717-61-3	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051789.D
 Acq On : 23 Dec 2021 17:44
 Operator : CG/JU
 Sample : M5215-04
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.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
Ethylbenzene	5.781	40.2	ng/ul	44576900	1	8.324	22183200	20.0
Bicyclo[2.1.1]h...	5.975	110.9	ng/ul	123051000	1	8.324	22183200	20.0
o-Xylene	6.304	55.9	ng/ul	61984900	1	8.324	22183200	20.0
(DEL) Alkane: S...	6.550	8.8	ng/ul	9715190	1	8.324	22183200	20.0
(DEL) Alkane: S...	6.668	4.2	ng/ul	4690990	1	8.324	22183200	20.0
Benzene, (1-met...	6.768	9.2	ng/ul	10191000	1	8.324	22183200	20.0
(DEL) Alkane: S...	6.844	5.6	ng/ul	6195780	1	8.324	22183200	20.0
(DEL) Alkane: C...	6.926	16.4	ng/ul	18220700	1	8.324	22183200	20.0
(DEL) Alkane: S...	7.191	4.7	ng/ul	5162160	1	8.324	22183200	20.0
Benzene, propyl-	7.273	15.1	ng/ul	16730700	1	8.324	22183200	20.0
(DEL) Alkane: S...	7.473	23.6	ng/ul	26172000	1	8.324	22183200	20.0
Benzene, 1,2,3-...	7.549	22.4	ng/ul	24866900	1	8.324	22183200	20.0
Benzene, 1-ethy...	7.708	13.2	ng/ul	14635600	1	8.324	22183200	20.0
Benzene, 1,2,4-...	7.984	59.5	ng/ul	65985100	1	8.324	22183200	20.0
Octane, 1,1'-ox...	8.413	6.2	ng/ul	6884660	1	8.324	22183200	20.0
Benzaldehyde, 4...	8.471	23.0	ng/ul	25546700	1	8.324	22183200	20.0
(DEL) Alkane: S...	8.530	5.7	ng/ul	6359030	1	8.324	22183200	20.0
1-Iodo-2-methyl...	8.589	9.5	ng/ul	10546400	1	8.324	22183200	20.0
(DEL) Alkane: C...	8.642	8.8	ng/ul	9779200	1	8.324	22183200	20.0
Indane	8.718	10.6	ng/ul	11791500	1	8.324	22183200	20.0
Benzene, 1-meth...	8.883	24.9	ng/ul	27654100	1	8.324	22183200	20.0
(DEL) Alkane: S...	8.935	5.0	ng/ul	5530300	1	8.324	22183200	20.0
(DEL) Alkane: S...	9.112	10.5	ng/ul	11639300	1	8.324	22183200	20.0
Naphthalene, de...	9.200	4.8	ng/ul	5315660	1	8.324	22183200	20.0
Benzene, 4-ethy...	9.288	3.0	ng/ul	3385380	1	8.324	22183200	20.0
p-Cymene	9.347	7.9	ng/ul	8759410	1	8.324	22183200	20.0
Benzene, 1-ethy...	9.540	7.4	ng/ul	8171540	1	8.324	22183200	20.0
(DEL) Alkane: S...	9.617	46.9	ng/ul	51968400	1	8.324	22183200	20.0
Benzene, 1-meth...	9.699	12.6	ng/ul	13937600	1	8.324	22183200	20.0
Benzene, 2-ethy...	9.793	6.5	ng/ul	5409230	2	11.138	16606700	20.0
unknown-01	9.822	7.2	ng/ul	5960180	2	11.138	16606700	20.0
(DEL) Alkane: S...	9.916	5.6	ng/ul	4612840	2	11.138	16606700	20.0
1,3,8-p-Menthat...	9.975	10.8	ng/ul	8981790	2	11.138	16606700	20.0
Benzene, 1,2,3,...	10.040	10.7	ng/ul	8884140	2	11.138	16606700	20.0
Naphthalene, de...	10.093	6.1	ng/ul	5100140	2	11.138	16606700	20.0
Benzene, 1,3-di...	10.334	9.5	ng/ul	7856050	2	11.138	16606700	20.0
(DEL) Alkane: S...	10.404	8.2	ng/ul	6793810	2	11.138	16606700	20.0
(DEL) Alkane: S...	10.475	6.5	ng/ul	5437020	2	11.138	16606700	20.0
Benzene, 1-meth...	10.551	14.3	ng/ul	11838100	2	11.138	16606700	20.0
3,5-Dimethylben...	10.604	3.5	ng/ul	2947370	2	11.138	16606700	20.0
(DEL) Alkane: S...	10.657	3.8	ng/ul	3110350	2	11.138	16606700	20.0
(DEL) Alkane: C...	11.832	2.2	ng/ul	1857150	2	11.138	16606700	20.0
(DEL) Alkane: S...	12.025	2.5	ng/ul	2074380	2	11.138	16606700	20.0
(DEL) Alkane: S...	12.131	4.6	ng/ul	3819000	2	11.138	16606700	20.0
(DEL) Alkane: S...	12.525	10.2	ng/ul	8492810	2	11.138	16606700	20.0
(DEL) Alkane: S...	13.130	13.3	ng/ul	690663	3	14.916	1036570	20.0
unknown-02	13.171	11.9	ng/ul	619273	3	14.916	1036570	20.0
(DEL) Alkane: S...	13.306	16.1	ng/ul	833779	3	14.916	1036570	20.0
(DEL) Alkane: S...	13.394	10.2	ng/ul	530638	3	14.916	1036570	20.0
(DEL) Alkane: S...	13.447	26.5	ng/ul	1373000	3	14.916	1036570	20.0
Naphthalene, 2,...	14.234	11.8	ng/ul	613330	3	14.916	1036570	20.0
unknown-03	15.709	10.2	ng/ul	527911	3	14.916	1036570	20.0

(DEL) Alkane: S...	16.631	20.0	ng/ul	623194	4	17.665	623194	20.0
unknown-04	31.499	19.7	ng/ul	502085	6	25.466	509287	20.0

Instrument :
BNA_G
ClientSampleId :
BGLD6