

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
 Data File : BG048198.D
 Acq On : 18 Feb 2021 16:44
 Operator : CG/JU
 Sample : M1429-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBGR3

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.251	321	325	332	rVB	45514	68275	0.49%	0.104%
2	5.656	388	394	400	rBV	28702	47526	0.34%	0.072%
3	6.196	477	486	499	rVB	678782	1194375	8.57%	1.811%
4	6.443	521	528	536	rBV4	23851	53488	0.38%	0.081%
5	6.590	544	553	559	rBV2	143394	260146	1.87%	0.394%
6	7.301	667	674	682	rBV4	58694	154400	1.11%	0.234%
7	7.372	682	686	690	rVV2	38729	67605	0.49%	0.103%
8	7.436	690	697	708	rVB	155289	350643	2.52%	0.532%
9	7.554	709	717	721	rBV5	74222	159269	1.14%	0.241%
10	7.607	721	726	730	rVV4	80475	158467	1.14%	0.240%
11	7.654	730	734	740	rVV	184716	322845	2.32%	0.490%
12	7.706	740	743	746	rVV3	57878	103533	0.74%	0.157%
13	7.748	747	750	755	rVV3	70239	139831	1.00%	0.212%
14	7.941	777	783	785	rBV	45289	77063	0.55%	0.117%
15	7.965	785	787	792	rVV3	36664	50926	0.37%	0.077%
16	8.112	806	812	818	rVB	182490	311400	2.23%	0.472%
17	8.852	931	938	944	rVV2	142231	264921	1.90%	0.402%
18	8.917	944	949	959	rVB8	28639	71279	0.51%	0.108%
19	9.040	964	970	976	rVB4	39230	69679	0.50%	0.106%
20	9.140	981	987	996	rVB3	68931	162076	1.16%	0.246%
21	9.258	996	1007	1011	rBV	728965	1369204	9.83%	2.076%
22	9.375	1019	1027	1033	rVB	562965	1024506	7.35%	1.553%
23	9.446	1033	1039	1044	rBV	236367	435352	3.12%	0.660%
24	9.528	1050	1053	1062	rVB6	35220	69748	0.50%	0.106%
25	9.628	1062	1070	1076	rBV6	88631	222582	1.60%	0.337%
26	9.692	1076	1081	1089	rVB	179734	299344	2.15%	0.454%
27	9.786	1089	1097	1101	rBV3	133321	284594	2.04%	0.432%
28	9.833	1101	1105	1107	rVV4	138857	219210	1.57%	0.332%
29	9.869	1107	1111	1124	rVB	357215	675370	4.85%	1.024%
30	10.010	1124	1135	1142	rBV6	274009	774310	5.56%	1.174%
31	10.198	1158	1167	1174	rBV	7702756	13934121	100.00%	21.128%
32	10.303	1177	1185	1192	rBV3	2575583	4973891	35.70%	7.542%
33	10.486	1211	1216	1219	rBV4	237904	497516	3.57%	0.754%
34	10.591	1224	1234	1243	rBV2	2332540	6548368	47.00%	9.929%

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35	10.738	1254	1259	1260	rBV	553027	697158	5.00%	1.057%
36	10.762	1260	1263	1270	rVB2	950947	1521275	10.92%	2.307%
37	10.838	1270	1276	1281	rBV	928576	1569469	11.26%	2.380%
38	10.885	1281	1284	1287	rVB3	141921	175482	1.26%	0.266%
39	10.932	1287	1292	1297	rBV3	417509	689683	4.95%	1.046%
40	10.985	1297	1301	1307	rVB	672248	1147016	8.23%	1.739%
41	11.044	1307	1311	1316	rBV2	97315	162849	1.17%	0.247%
42	11.097	1316	1320	1326	rVB2	228818	381017	2.73%	0.578%
43	11.208	1336	1339	1348	rVB3	91535	156878	1.13%	0.238%
44	11.314	1350	1357	1367	rBV3	156281	334582	2.40%	0.507%
45	11.402	1369	1372	1378	rVB3	36504	59495	0.43%	0.090%
46	11.484	1379	1386	1397	rBV3	89358	275568	1.98%	0.418%
47	11.996	1467	1473	1479	rBV2	153784	309770	2.22%	0.470%
48	12.054	1479	1483	1487	rVV3	60846	92143	0.66%	0.140%
49	12.283	1515	1522	1532	rBV4	165881	384935	2.76%	0.584%
50	12.372	1532	1537	1545	rVB2	126422	220333	1.58%	0.334%
51	12.466	1548	1553	1556	rBV	79660	121922	0.87%	0.185%
52	12.560	1559	1569	1574	rVB	332670	700984	5.03%	1.063%
53	12.665	1577	1587	1589	rBV3	157287	382253	2.74%	0.580%
54	12.795	1600	1609	1613	rBV5	162588	397386	2.85%	0.603%
55	12.836	1613	1616	1623	rVB	155211	265594	1.91%	0.403%
56	12.912	1623	1629	1632	rBV2	177083	301533	2.16%	0.457%
57	12.959	1632	1637	1642	rVV	776760	1212835	8.70%	1.839%
58	13.012	1642	1646	1649	rVV	271615	417631	3.00%	0.633%
59	13.041	1649	1651	1658	rVB4	129874	177355	1.27%	0.269%
60	13.153	1663	1670	1678	rVV3	201763	412464	2.96%	0.625%
61	13.500	1724	1729	1733	rBV6	31181	58325	0.42%	0.088%
62	13.582	1738	1743	1748	rVV3	91779	196717	1.41%	0.298%
63	13.635	1748	1752	1760	rVV3	91636	210153	1.51%	0.319%
64	13.705	1760	1764	1772	rVV4	73591	199787	1.43%	0.303%
65	13.793	1772	1779	1789	rVB	2160975	3239716	23.25%	4.912%
66	13.923	1795	1801	1808	rBV5	108909	262483	1.88%	0.398%
67	14.017	1813	1817	1821	rVV2	124613	215342	1.55%	0.327%
68	14.064	1821	1825	1833	rVV5	90586	191026	1.37%	0.290%
69	14.146	1833	1839	1844	rVV	520789	765837	5.50%	1.161%
70	14.205	1844	1849	1855	rVB3	158850	249846	1.79%	0.379%
71	14.293	1860	1864	1870	rBV4	46315	93104	0.67%	0.141%

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72	14.387	1876	1880	1883	rBV4	31717	49972	0.36%	0.076%
73	14.440	1883	1889	1894	rVB	528543	834642	5.99%	1.266%
74	14.510	1895	1901	1905	rBV5	74307	131911	0.95%	0.200%
75	14.675	1926	1929	1934	rVB5	32835	50084	0.36%	0.076%
76	14.739	1934	1940	1946	rVV	381179	589773	4.23%	0.894%
77	14.804	1946	1951	1956	rVB	136935	197239	1.42%	0.299%
78	14.945	1971	1975	1978	rVV3	50191	81500	0.58%	0.124%
79	14.980	1978	1981	1984	rVV2	64564	107251	0.77%	0.163%
80	15.033	1984	1990	2001	rVB	553420	1007248	7.23%	1.527%
81	15.491	2063	2068	2072	rBV5	24888	50260	0.36%	0.076%
82	15.732	2103	2109	2117	rBV	717797	1140232	8.18%	1.729%
83	15.856	2124	2130	2137	rVB	281346	450320	3.23%	0.683%
84	16.244	2192	2196	2200	rBV	42413	58903	0.42%	0.089%
85	16.361	2212	2216	2218	rBV3	40810	51226	0.37%	0.078%
86	16.461	2229	2233	2236	rBV2	52466	71152	0.51%	0.108%
87	16.502	2237	2240	2245	rVB3	54931	73968	0.53%	0.112%
88	17.048	2328	2333	2341	rVB	181937	264117	1.90%	0.400%
89	17.307	2372	2377	2384	rVB	74907	139249	1.00%	0.211%
90	17.483	2402	2407	2413	rVB	461494	685865	4.92%	1.040%
91	17.583	2418	2424	2431	rVB2	802685	1212191	8.70%	1.838%
92	17.701	2440	2444	2451	rBV4	44443	86642	0.62%	0.131%
93	18.147	2517	2520	2525	rVB4	50335	75024	0.54%	0.114%
94	18.365	2554	2557	2563	rBV2	81721	127886	0.92%	0.194%
95	18.453	2568	2572	2578	rVB2	96931	146801	1.05%	0.223%
96	19.052	2670	2674	2679	rBV6	43282	92249	0.66%	0.140%
97	19.863	2807	2812	2818	rVB	1003266	1445855	10.38%	2.192%
98	21.761	3130	3135	3141	rBV2	484334	801844	5.75%	1.216%
99	24.822	3647	3656	3665	rVB	498683	1402622	10.07%	2.127%
100	25.045	3685	3694	3703	rVB	295250	859650	6.17%	1.303%

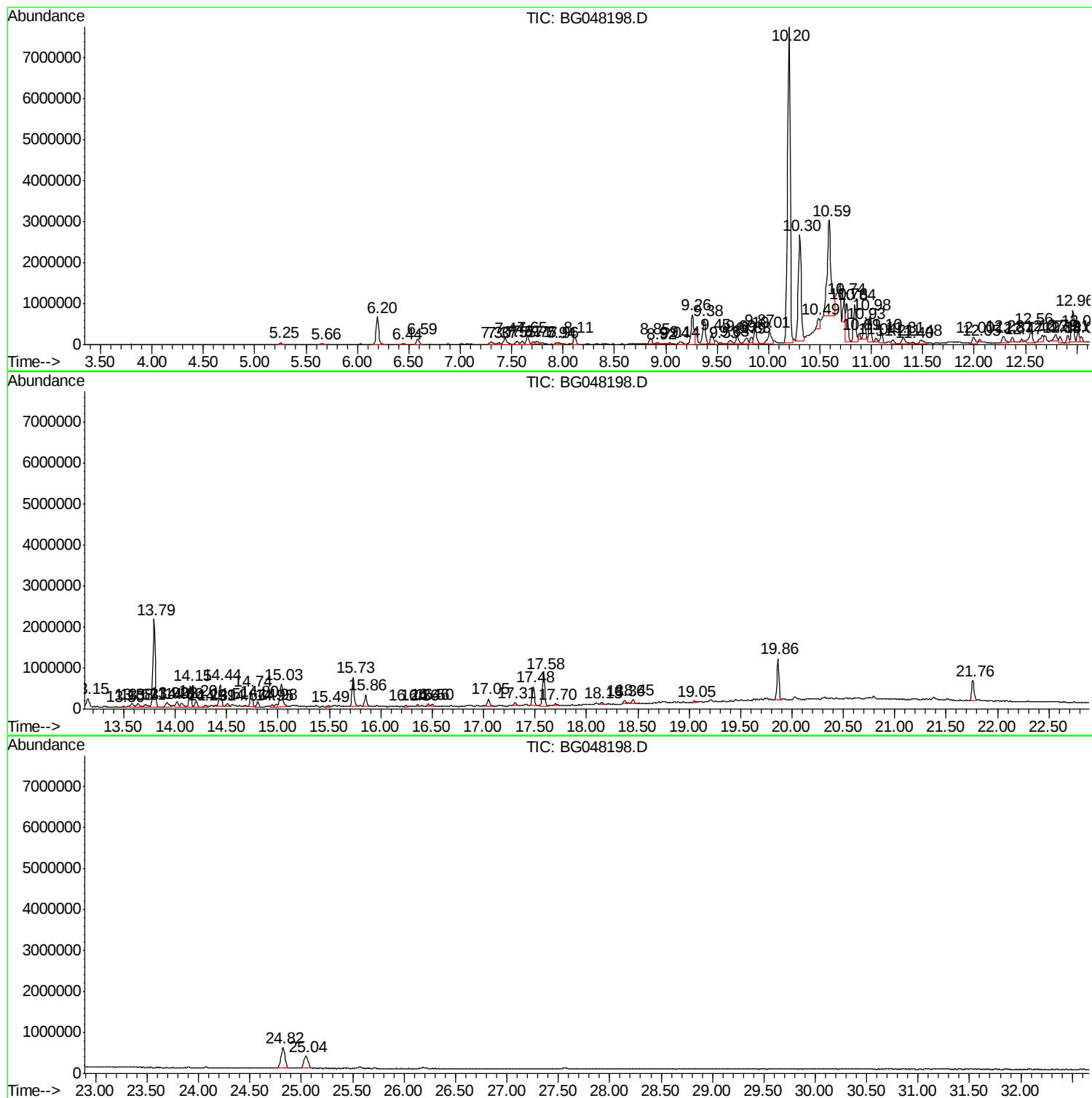
Sum of corrected areas: 65951485

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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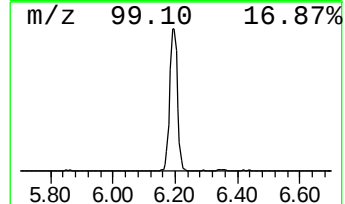
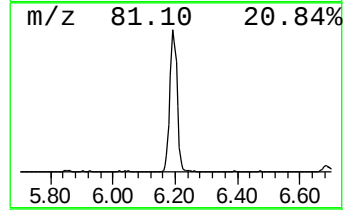
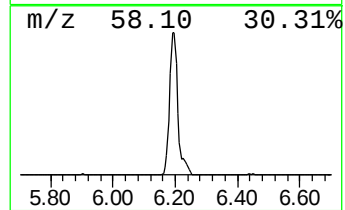
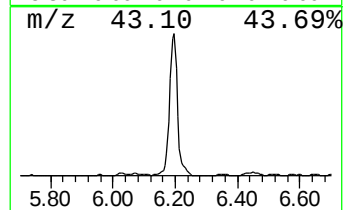
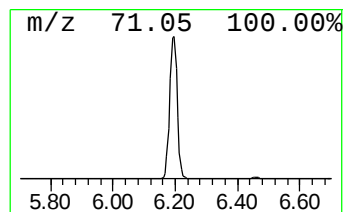
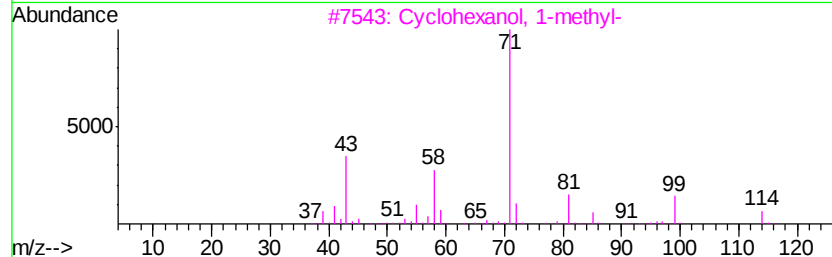
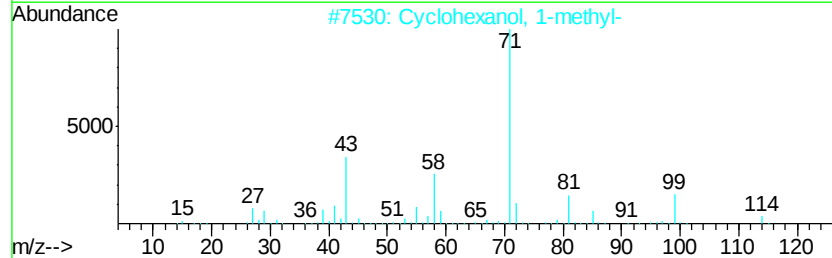
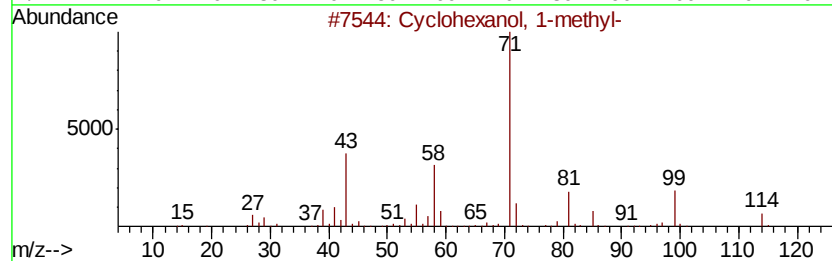
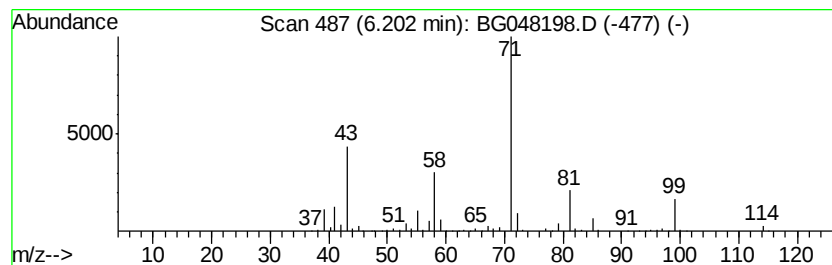
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexanol, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	76.71 ng/ul	1194380	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	94
2			Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	91
3			Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	78
4			2,4-Dimethyl-1-penten-3-ol	114	C7H14O	019781-54-5	58
5			trans-2-Methyl-4-hexen-3-ol	114	C7H14O	096346-76-8	53



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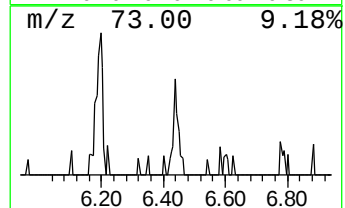
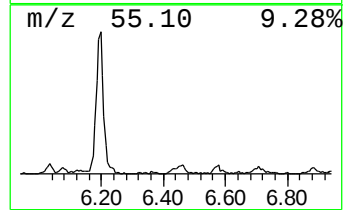
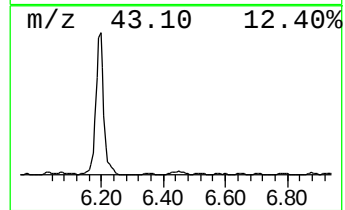
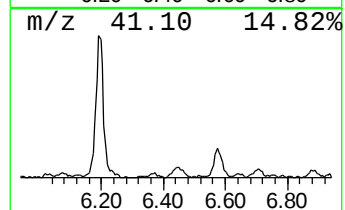
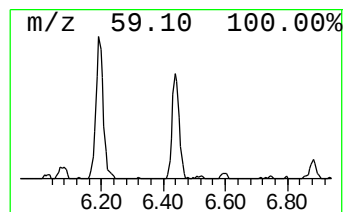
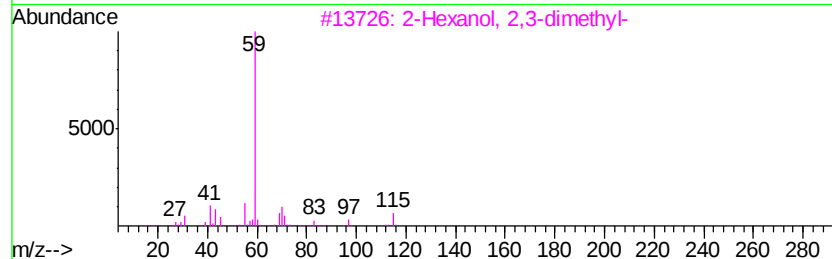
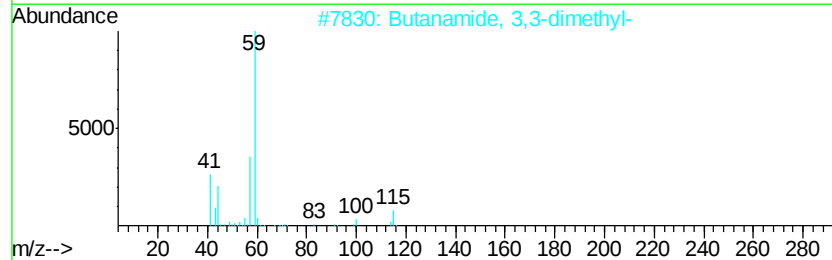
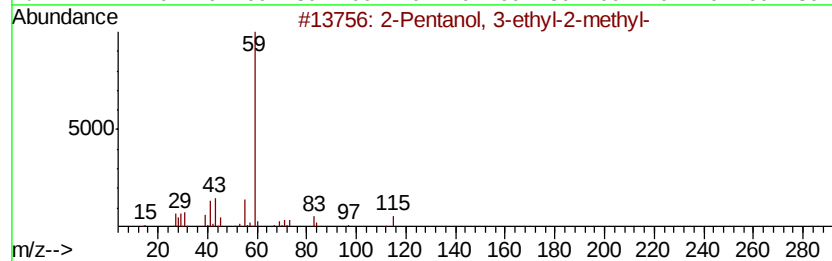
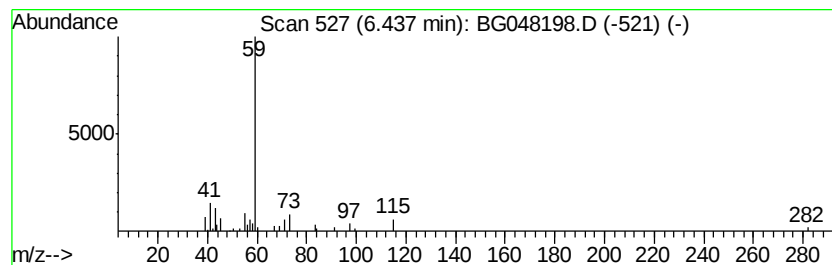
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanol, 3-ethyl-2-methyl- Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	3.44 ng/ul	53488	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanol, 3-ethyl-2-methyl-	130	C8H18O	019780-63-3	78
2		Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	53
3		2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	50
4		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	42



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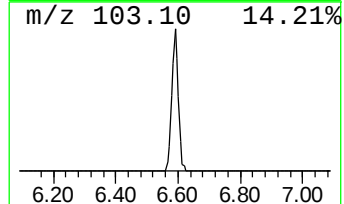
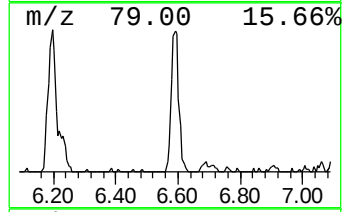
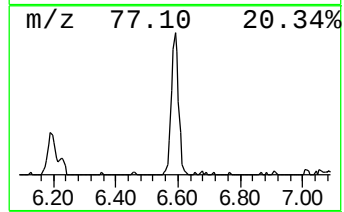
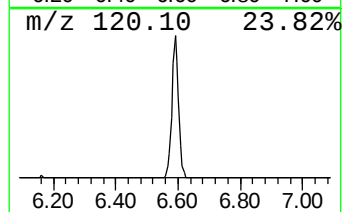
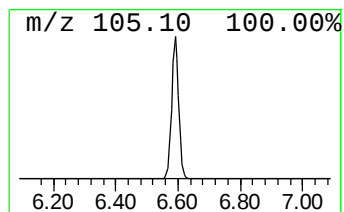
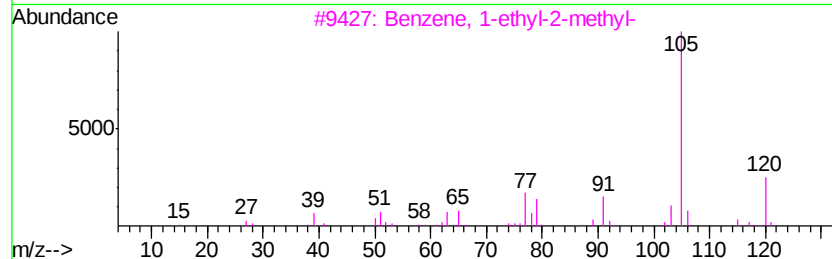
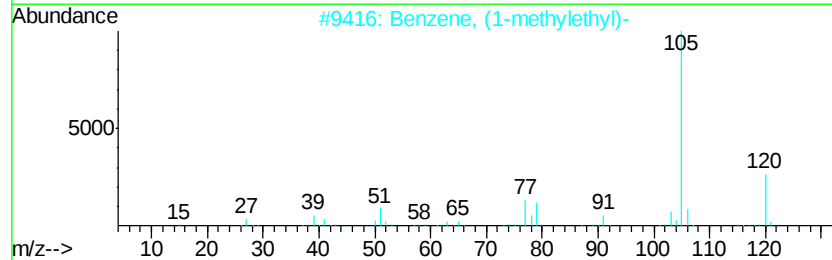
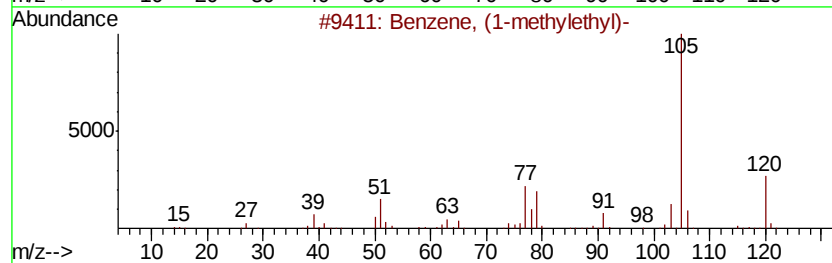
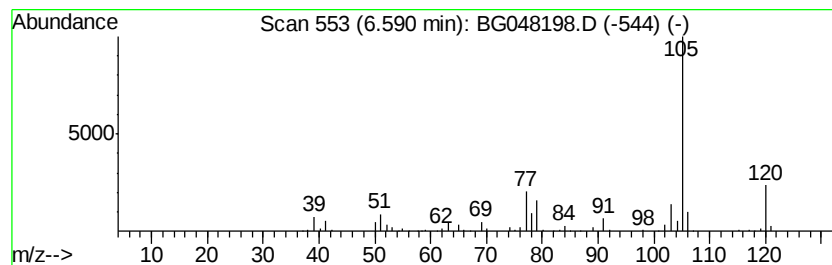
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, (1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	16.71 ng/ul	260146	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	83
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	80



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Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

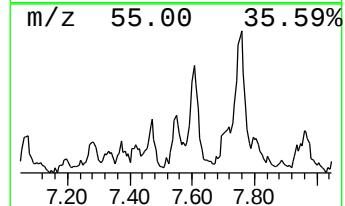
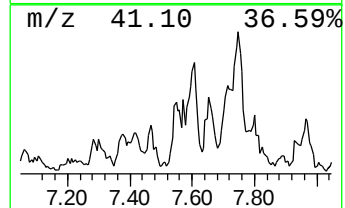
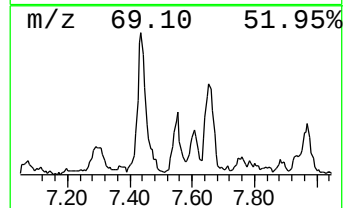
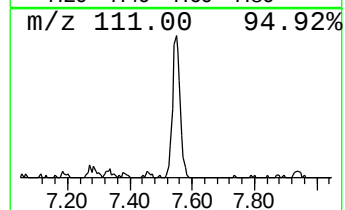
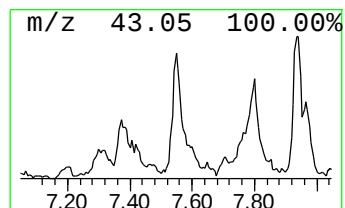
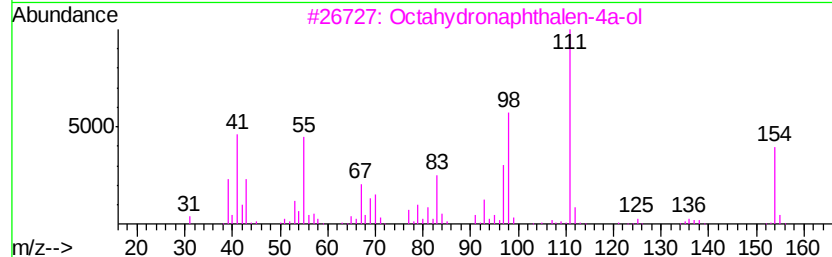
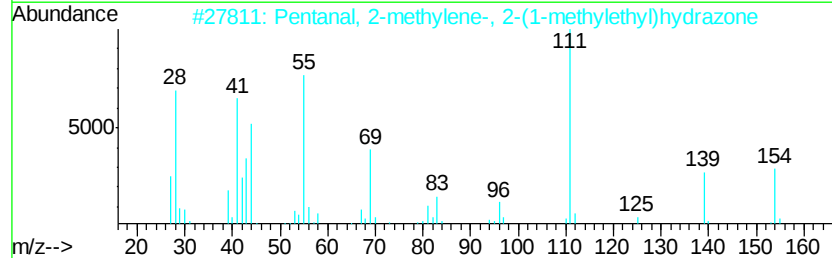
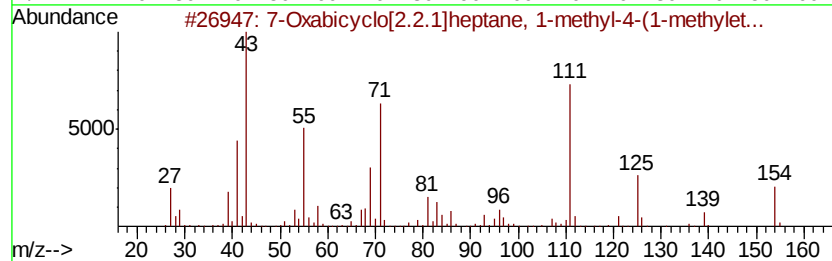
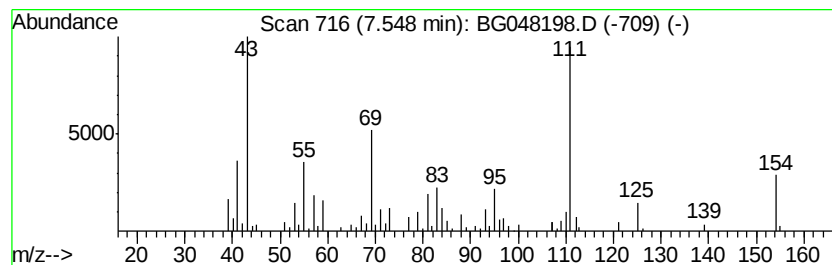
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.		
7.55	10.23 ng/ul	159269	1,4-Dichlorobenzene-d4	8.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Oxabicyclo[2.2.1]heptane, 1-me...	154	C10H18O	000470-67-7	46	
2	Pentanal, 2-methylene-, 2-(1-met...	154	C9H18N2	033063-80-8	43	
3	Octahydronaphthalen-4a-ol	154	C10H18O	055693-34-0	43	
4	4a(2H)-Naphthalenol, octahydro-, ...	154	C10H18O	001654-87-1	38	
5	Cyanamide, dibutyl-	154	C9H18N2	002050-54-6	38	



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Sample : M1429-03
Misc :
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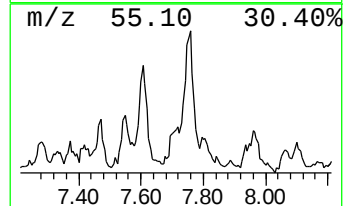
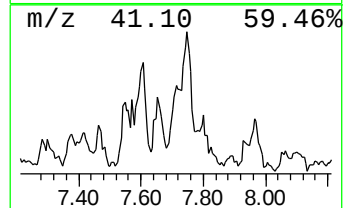
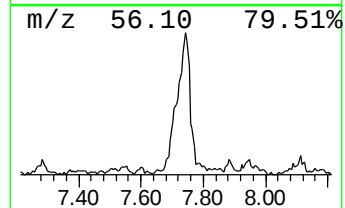
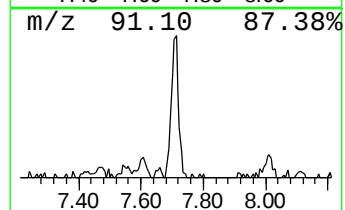
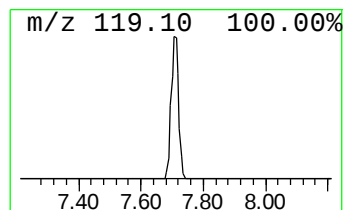
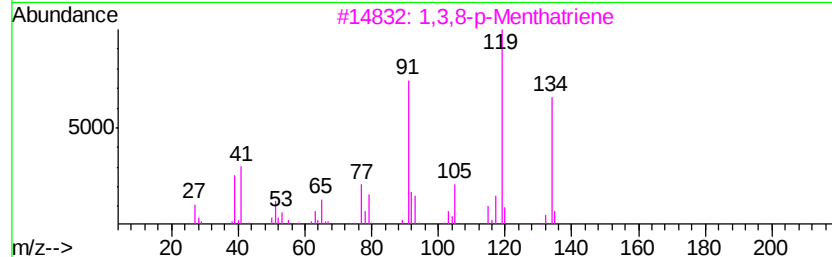
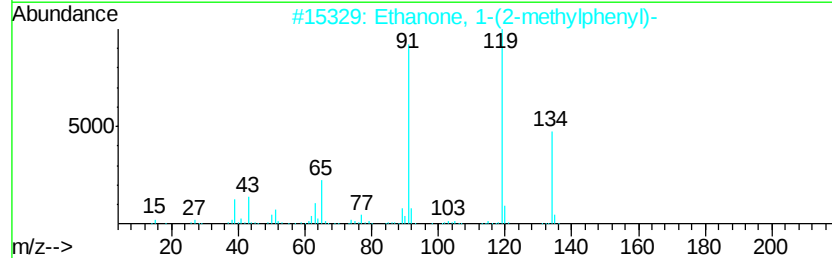
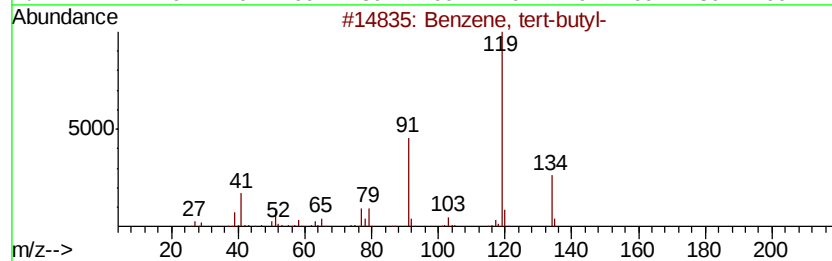
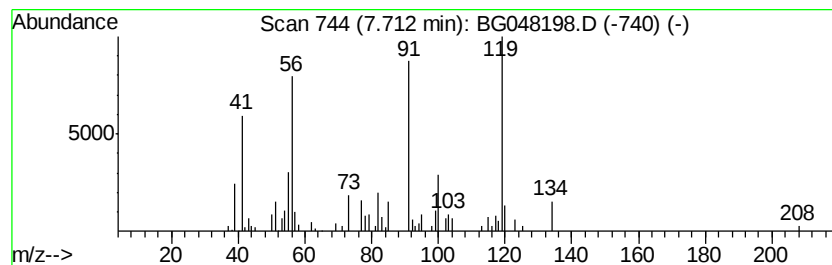
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown-02 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.71	6.65 ng/ul	103533	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, tert-butyl-	134	C10H14	000098-06-6	43
2		Ethanone, 1-(2-methylphenyl)-	134	C9H10O	000577-16-2	43
3		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	43
4		Ethanone, 1-(3-methylphenyl)-	134	C9H10O	000585-74-0	43
5		p-Cymene	134	C10H14	000099-87-6	38



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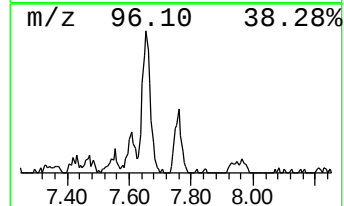
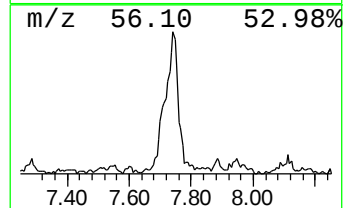
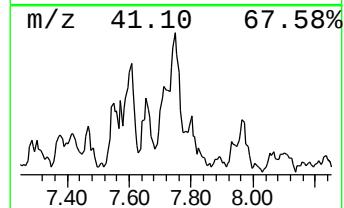
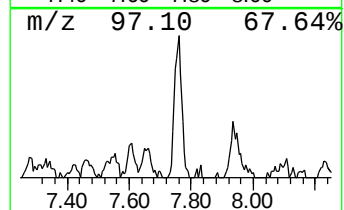
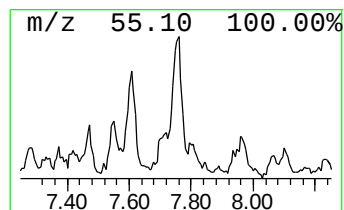
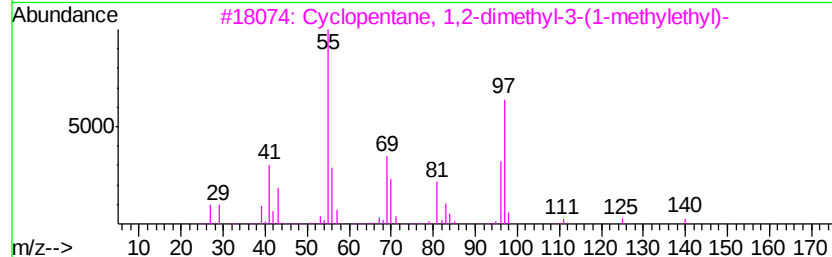
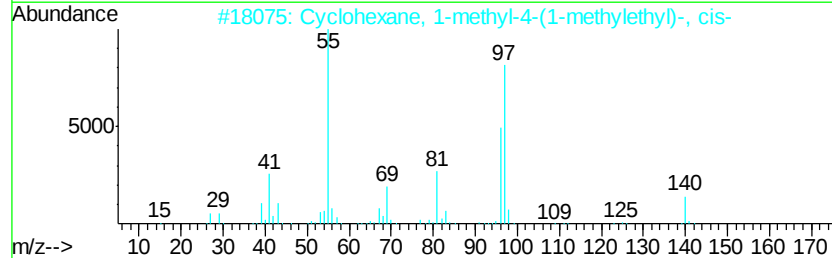
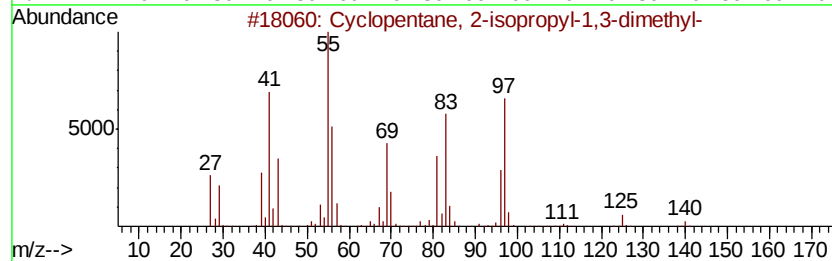
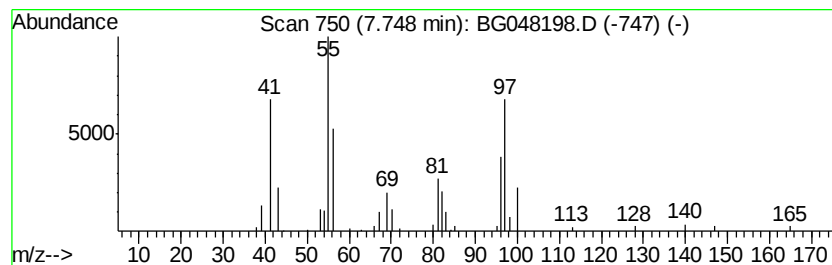
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Cyclic7.75 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	8.98 ng/ul	139831	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	62
2			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	53
3			Cyclopentane, 1,2-dimethyl-3-(1-...	140	C10H20	000489-20-3	52
4			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	47
5			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
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Misc :
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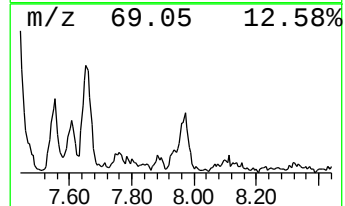
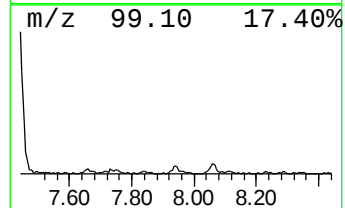
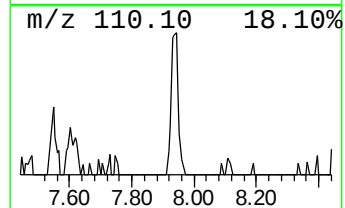
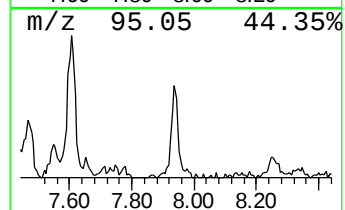
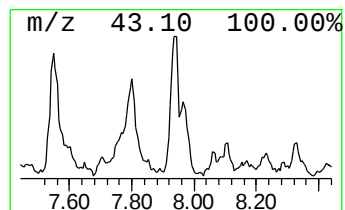
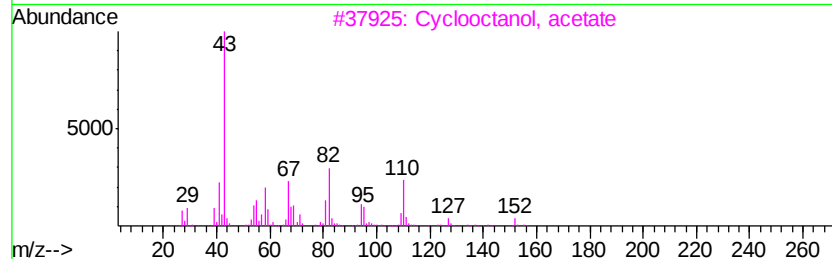
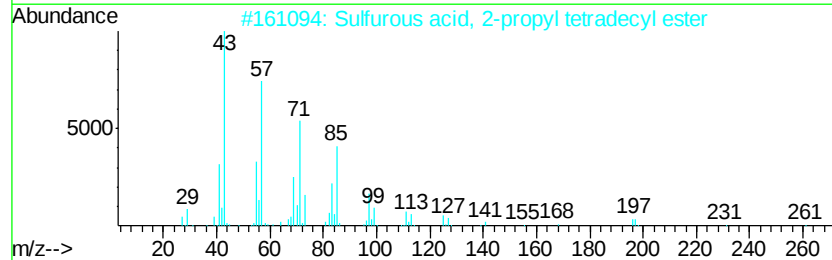
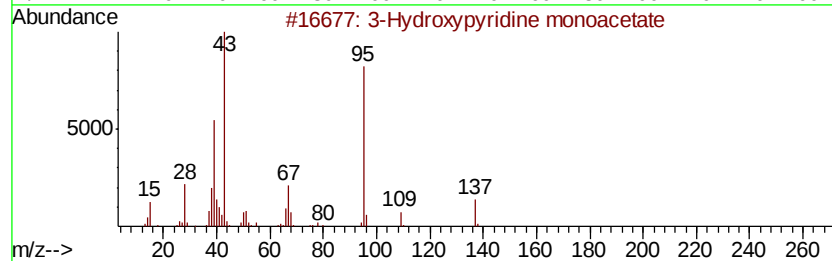
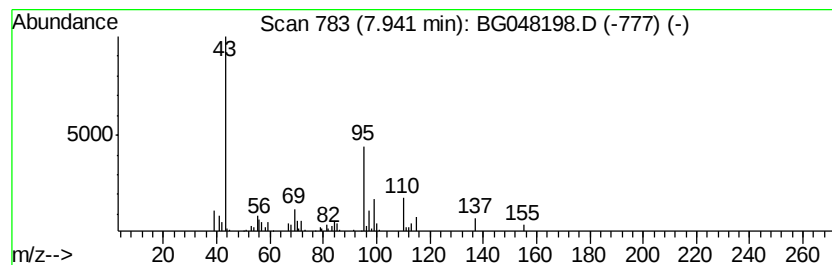
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown-03 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	4.95 ng/ul	77063	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Hydroxypyrindine monoacetate	137	C7H7NO2	017747-43-2	23
2			Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	9
3			Cyclooctanol, acetate	170	C10H18O2	000772-60-1	9
4			2-Oxabicyclo[3.2.1]nonan-7-one, ...	154	C9H14O2	013747-98-3	9
5			3-Octene-2,6-dione, 5,5,7-trimet...	182	C11H18O2	077142-73-5	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
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Misc :
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ClientSampleId :
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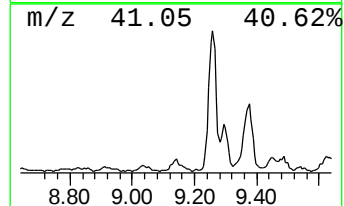
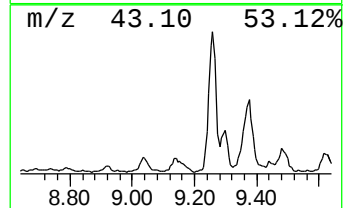
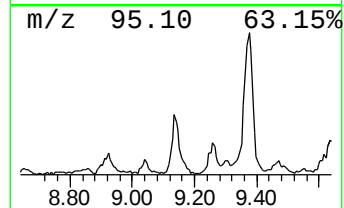
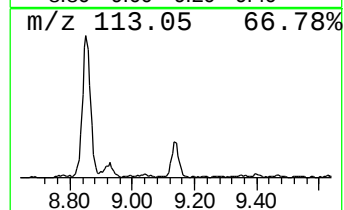
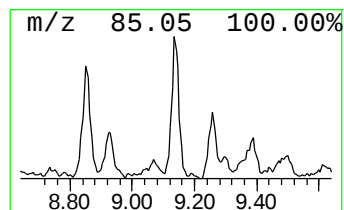
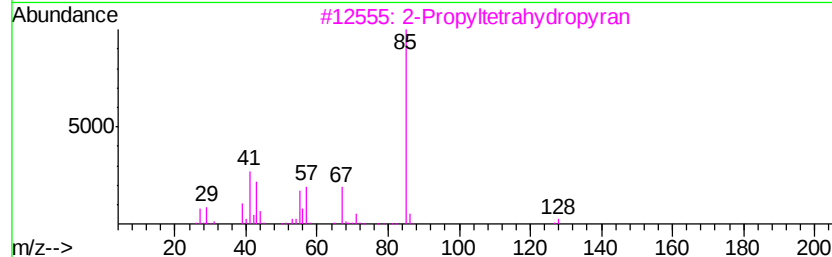
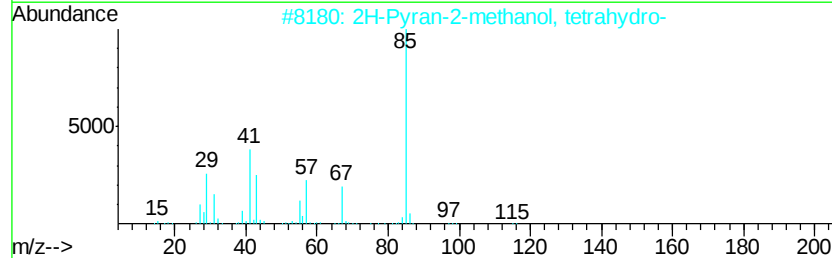
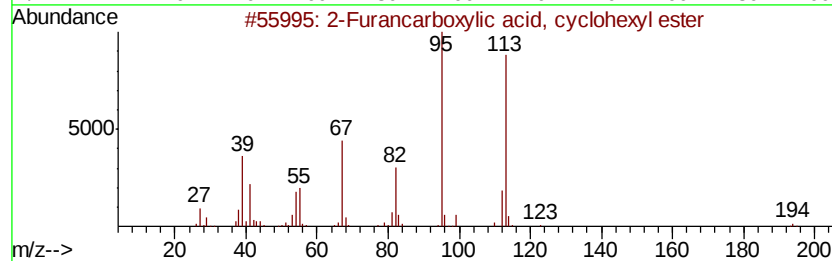
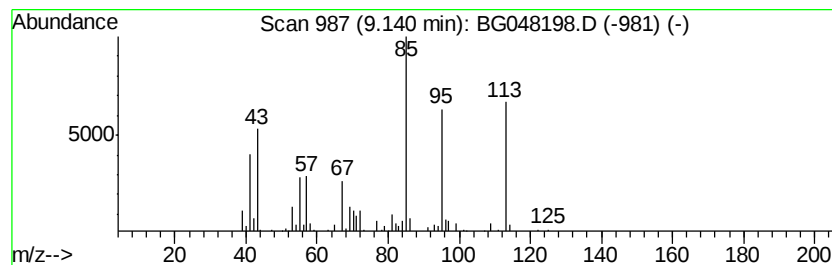
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown-04 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.14	10.41 ng/ul	162076	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Furancarboxylic acid, cyclohex...	194	C11H14O3	039251-91-7	38
2		2H-Pyran-2-methanol, tetrahydro-	116	C6H12O2	000100-72-1	35
3		2-Propyltetrahydropyran	128	C8H16O	003857-17-8	35
4		5,6-Epoxyhexanol-1	116	C6H12O2	021915-57-1	35
5		2-Furancarboxylic acid, decyl ester	252	C15H24O3	039251-90-6	35



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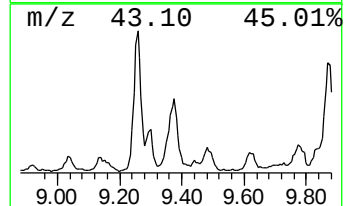
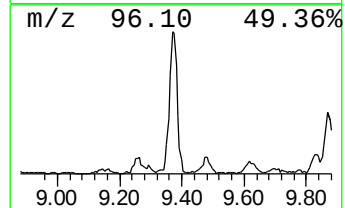
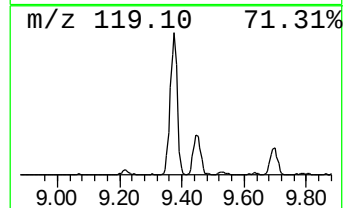
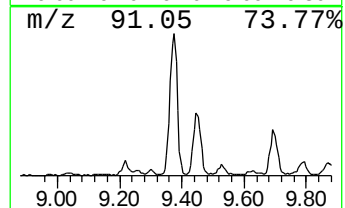
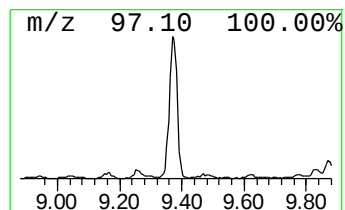
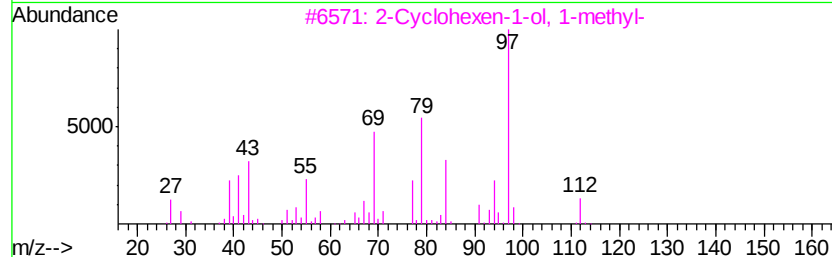
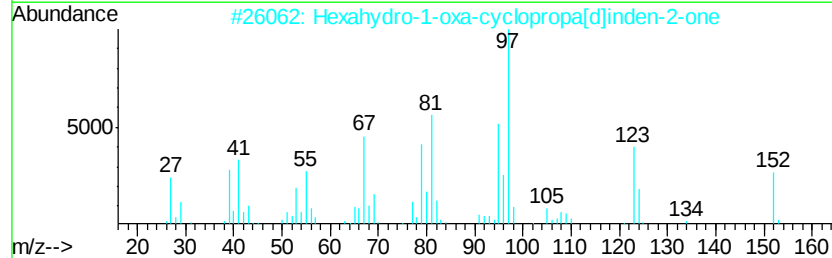
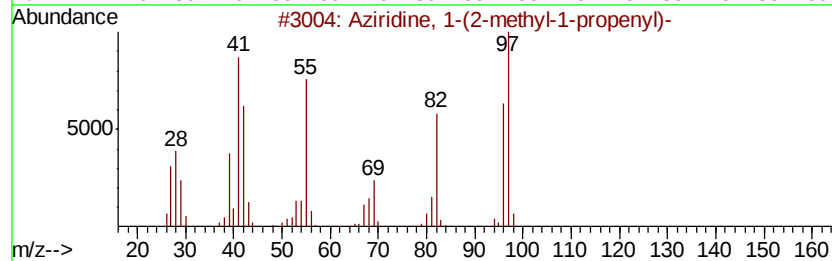
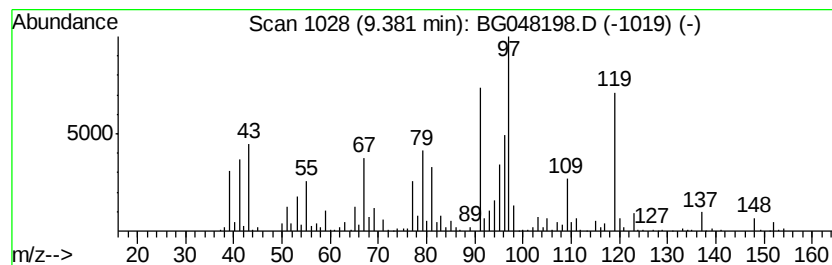
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown-05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	65.80 ng/ul	1024510	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aziridine, 1-(2-methyl-1-propenyl)-	97	C6H11N	080839-93-6	35
2		Hexahydro-1-oxa-cyclopropa[d]ind...	152	C9H12O2	037643-23-5	32
3		2-Cyclohexen-1-ol, 1-methyl-	112	C7H12O	023758-27-2	32
4		Glycine, furan-2-yl-methyl ester	155	C7H9NO3	1000306-78-7	32
5		m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
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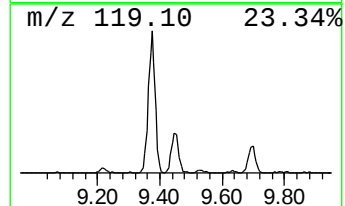
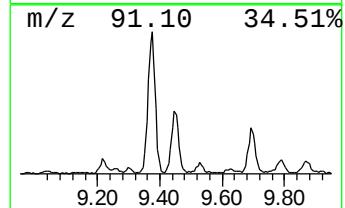
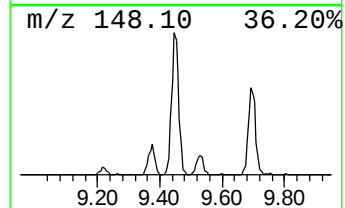
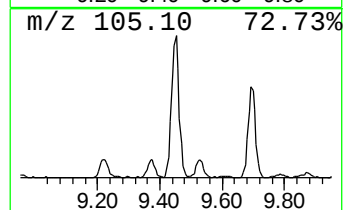
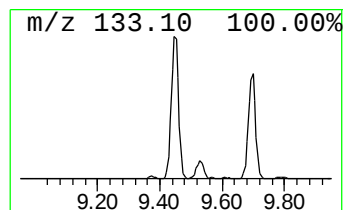
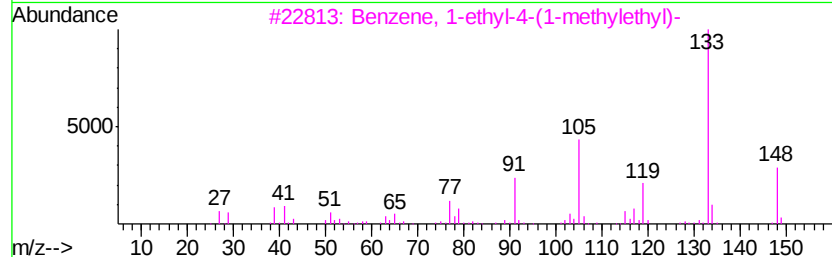
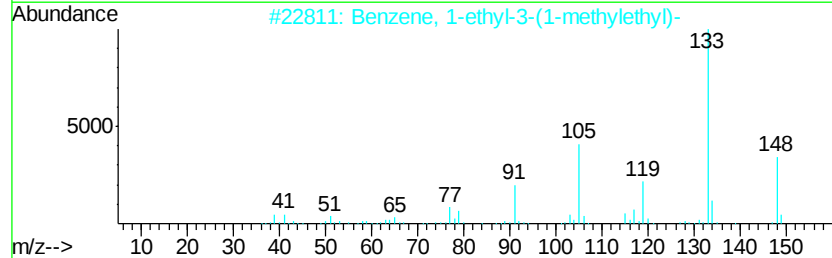
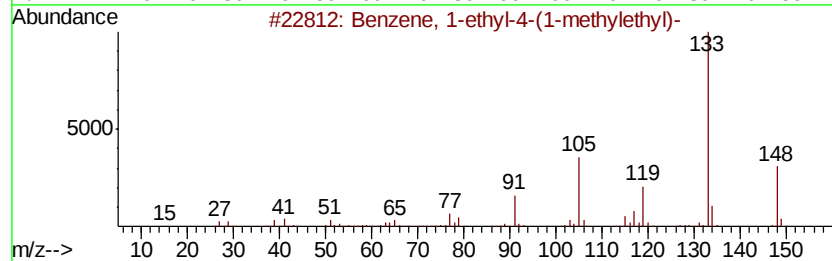
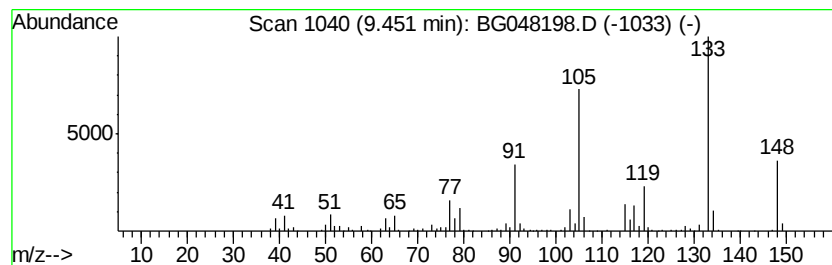
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.45	27.96 ng/ul	435352	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	97
2		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	96
3		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	94
4		Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	92
5		Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBG3

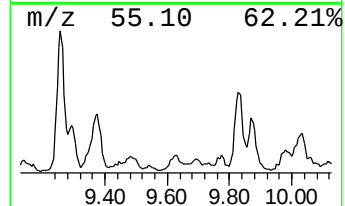
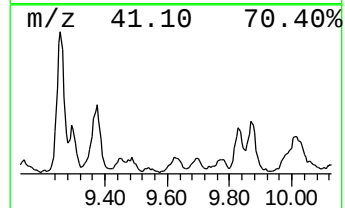
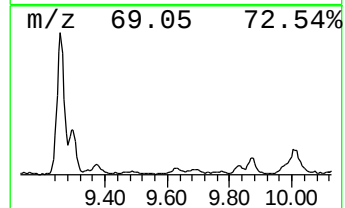
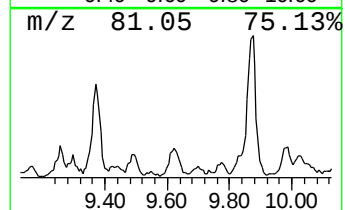
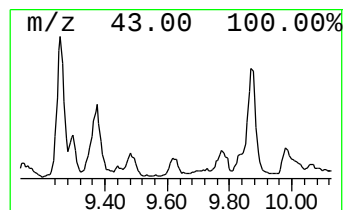
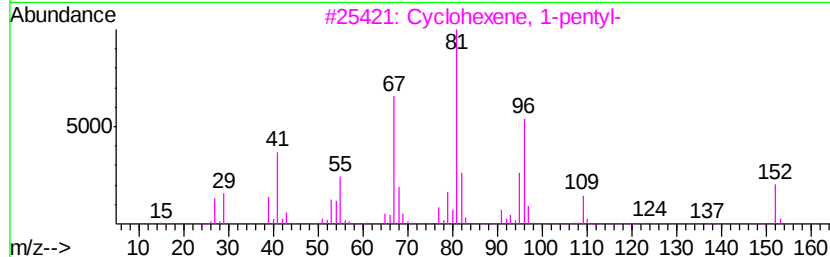
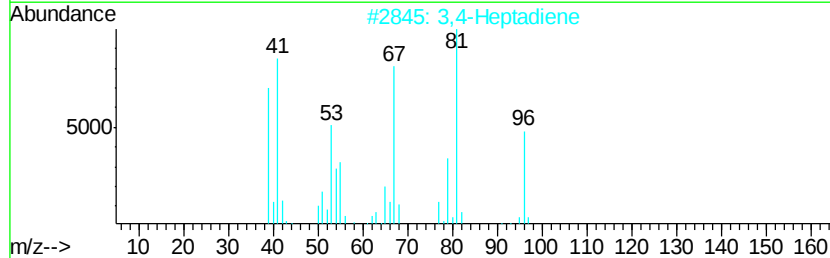
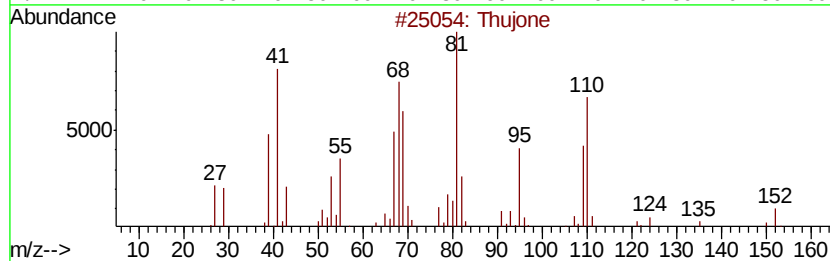
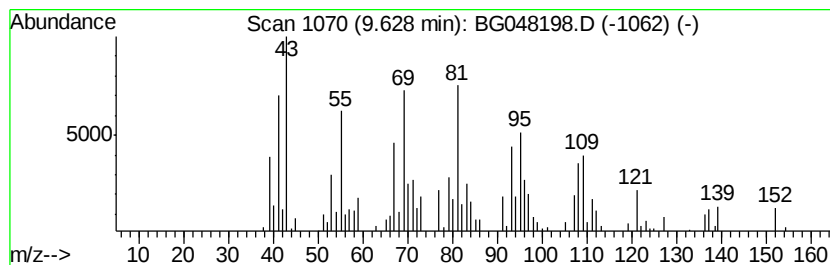
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 unknown-06 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	6.45 ng/ul	222582	Naphthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thujone	152	C10H16O	000546-80-5	43
2			3,4-Heptadiene	96	C7H12	002454-31-1	38
3			Cyclohexene, 1-pentyl-	152	C11H20	015232-85-6	38
4			Cyclohexaneethanol, 4-methyl-.be...	154	C10H18O	015714-12-2	38
5			Bicyclo[2.2.1]heptan-2-one, 4,7,...	152	C10H16O	010292-98-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

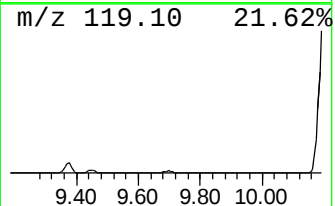
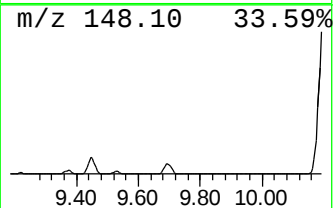
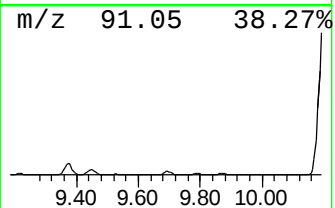
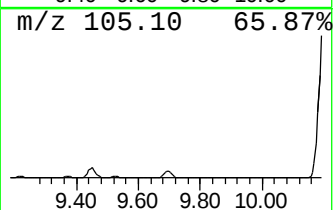
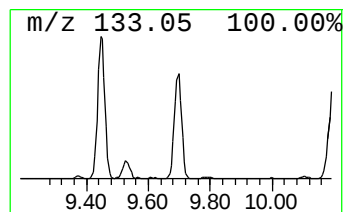
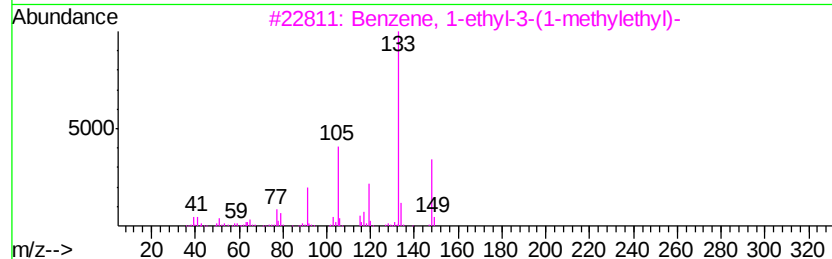
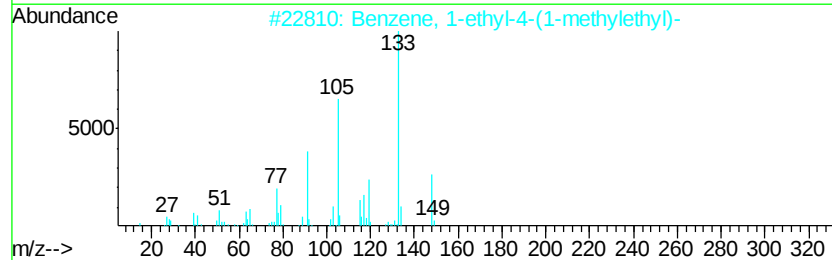
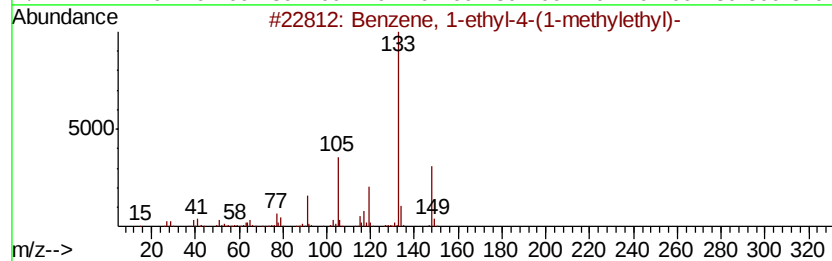
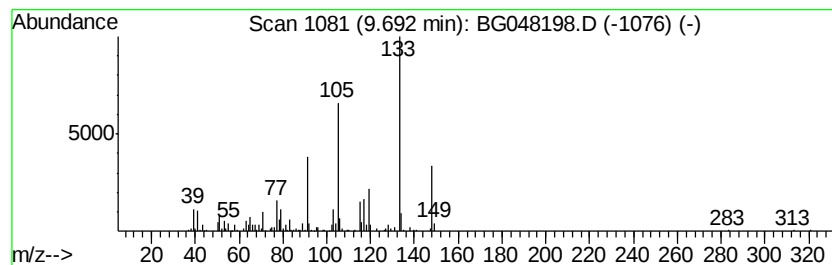
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	8.68 ng/ul	299344	Naphthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	97
2			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	96
3			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	94
4			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	91
5			Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

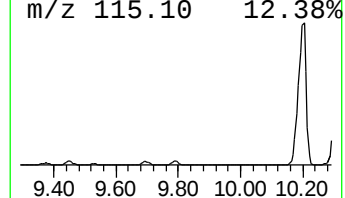
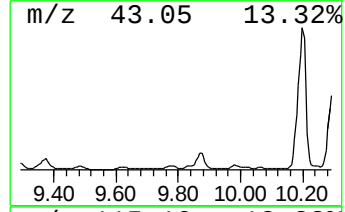
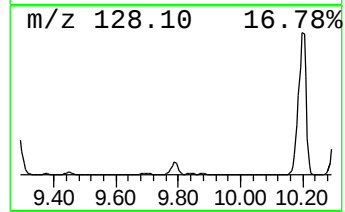
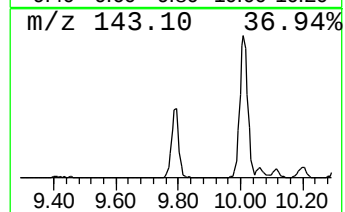
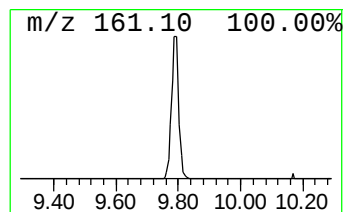
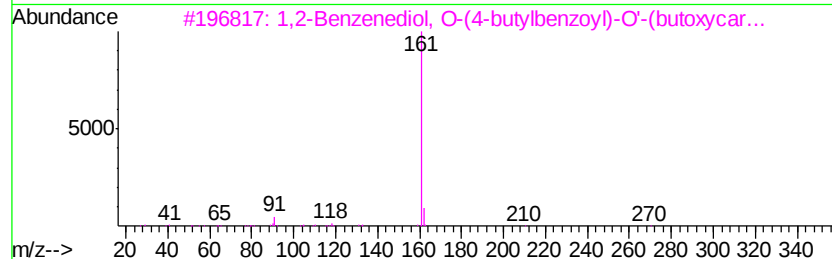
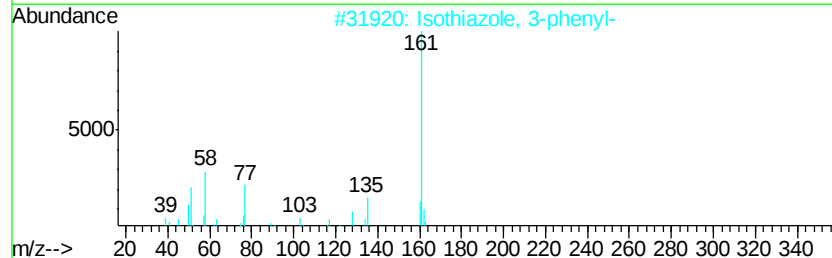
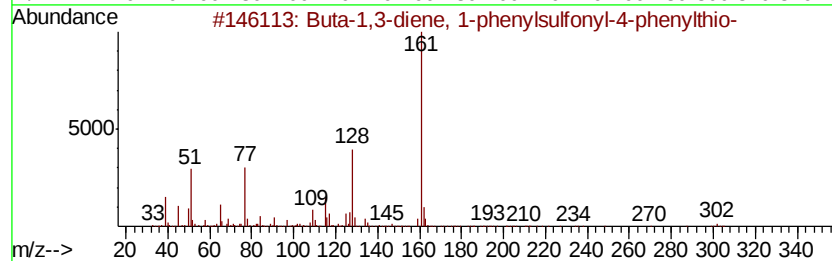
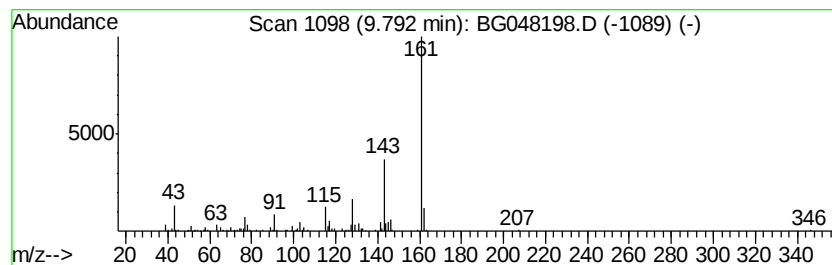
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Buta-1,3-diene, 1-phenylsul... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	8.25 ng/ul	284594	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Buta-1,3-diene, 1-phenylsulfonyl...	302	C16H14O2S2	121997-11-3	53
2		Isothiazole, 3-phenyl-	161	C9H7NS	010514-34-8	50
3		1,2-Benzenediol, O-(4-butylbenzo...	370	C22H26O5	1000329-73-7	47
4		1,3-Benzenediol, o-(4-butylbenzo...	364	C22H20O5	1000330-81-5	47
5		Phenylcyanide, 4-methoxy-2,6-dim...	161	C10H11NO	019111-77-4	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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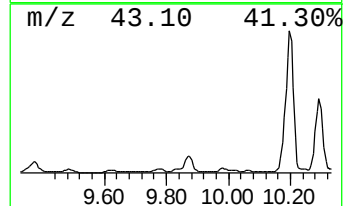
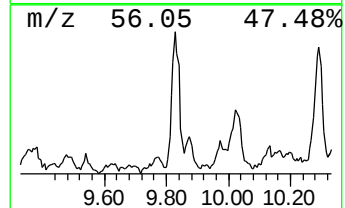
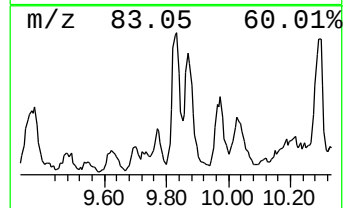
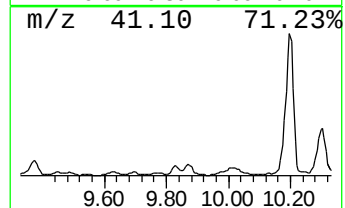
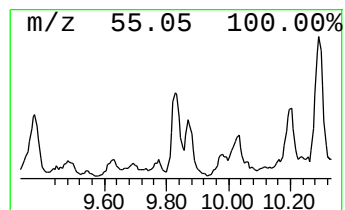
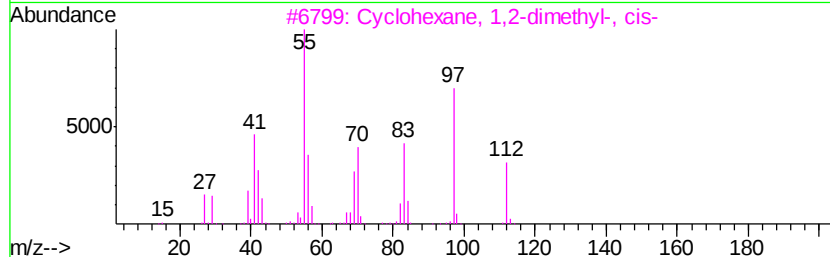
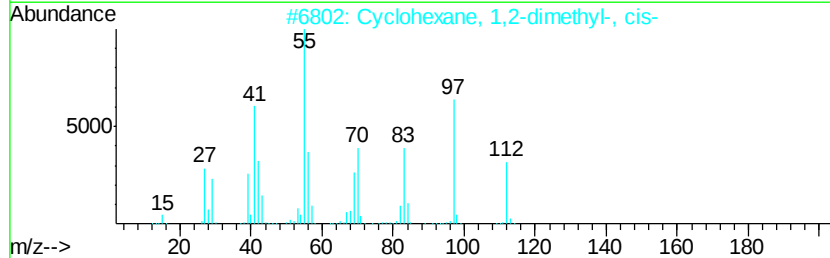
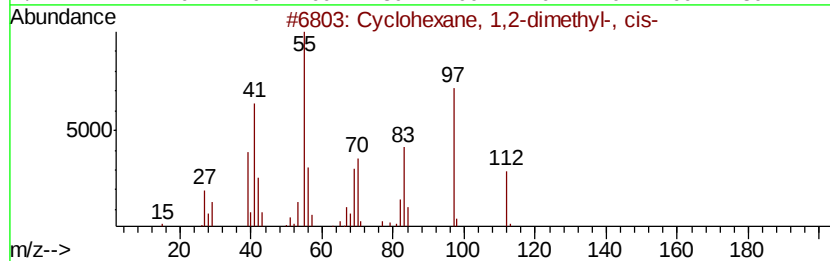
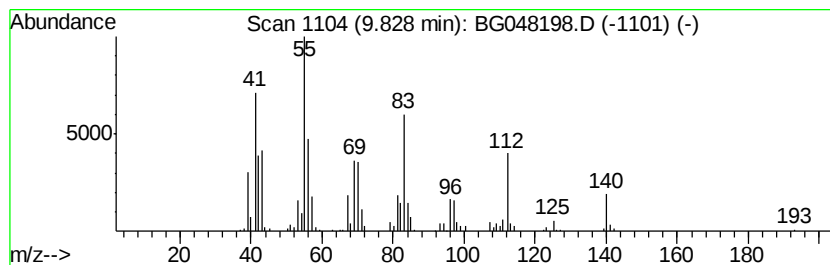
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Cyclic9.83 Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	6.36 ng/ul	219210	Napthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	58
2			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	52
3			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	52
4			1,2-Cyclohexanedione	112	C6H8O2	000765-87-7	49
5			4-Propylcyclohexanone	140	C9H16O	040649-36-3	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

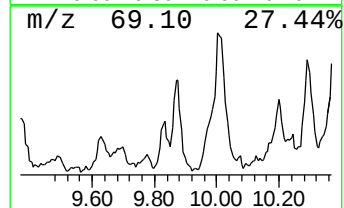
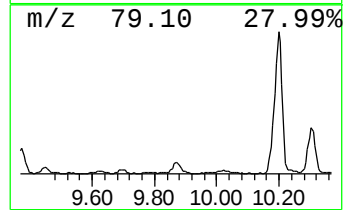
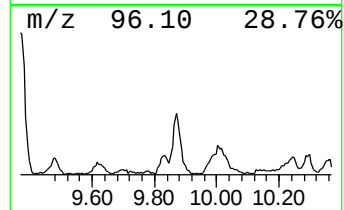
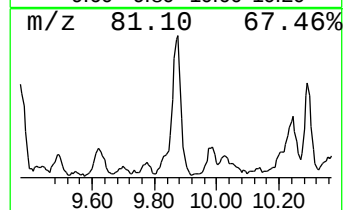
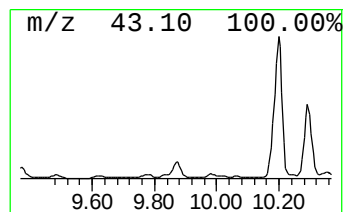
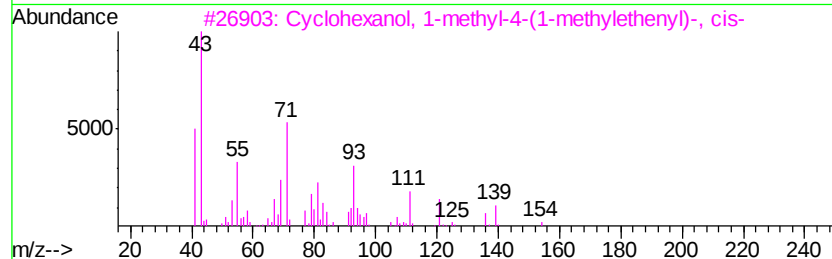
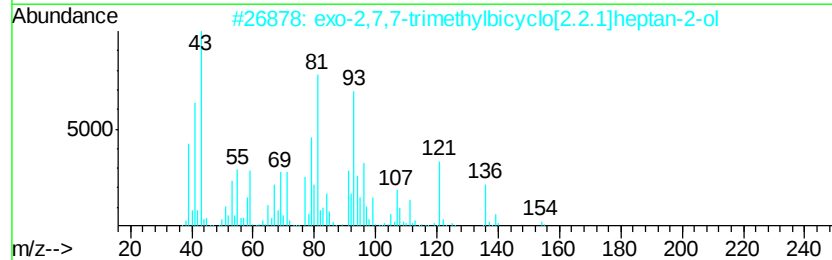
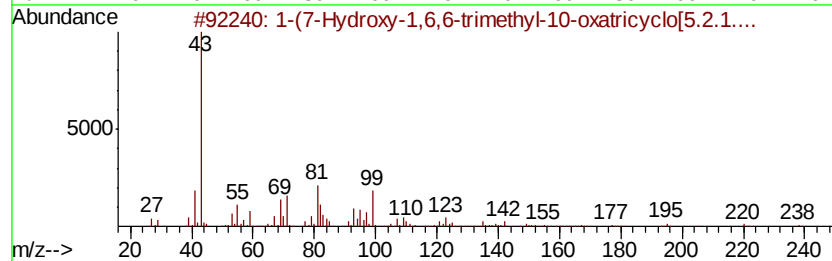
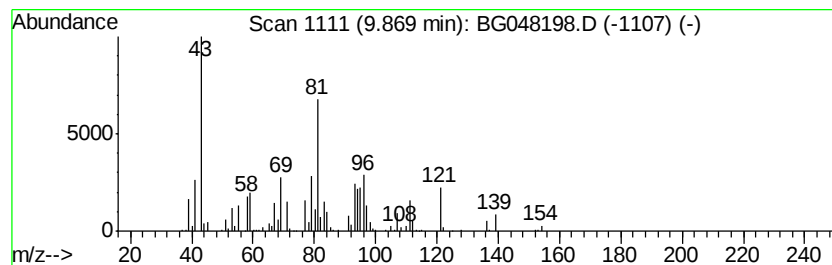
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 unknown-07 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	19.58 ng/ul	675370	Napthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(7-Hydroxy-1,6,6-trimethyl-10-...	238	C14H22O3	090165-01-8	47
2		exo-2,7,7-trimethylbicyclo[2.2.1]...	154	C10H18O	1000374-19-3	38
3		Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	007299-41-4	35
4		3-Cyclohexen-1-ol, 1-methyl-4-(1...	154	C10H18O	000586-82-3	27
5		2,6-Dimethyl-1,6-heptadien-4-ol ...	182	C11H18O2	070187-91-6	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

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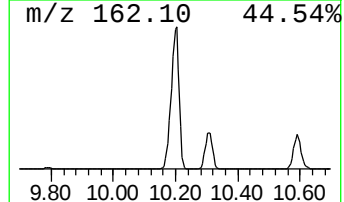
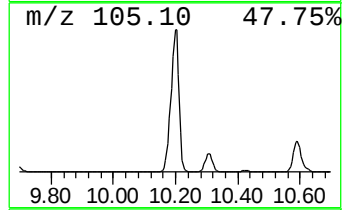
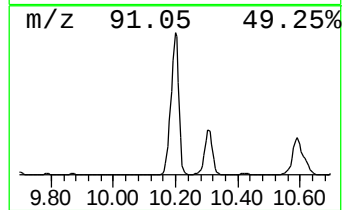
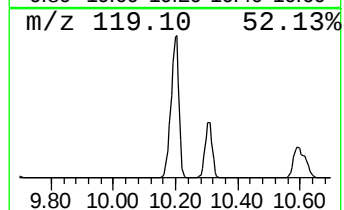
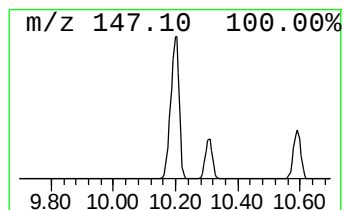
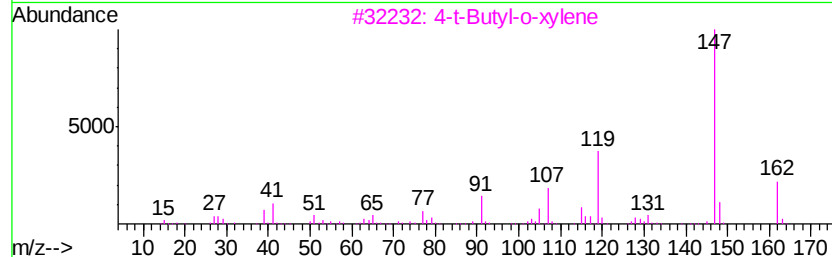
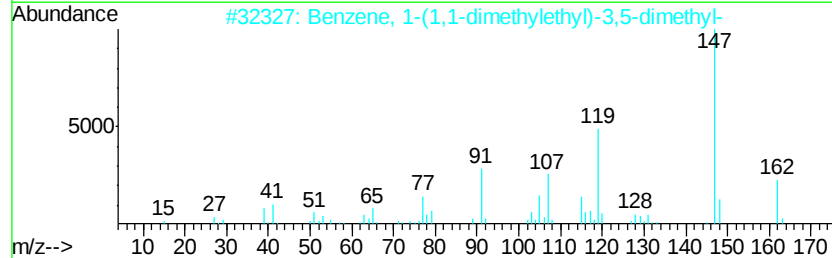
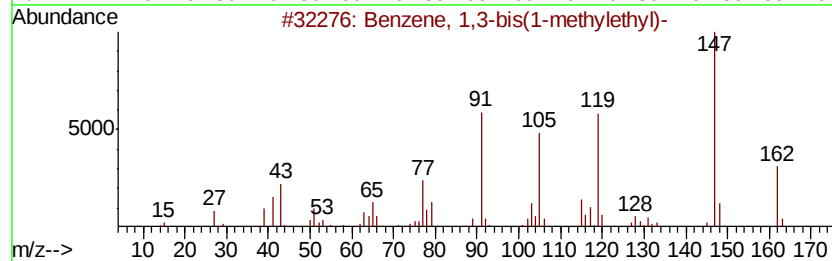
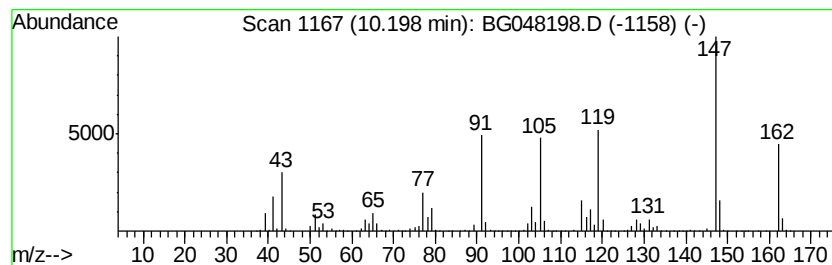
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Benzene, 1,3-bis(1-methylet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.20	404.07 ng/ul	13934100	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	95
2		Benzene, 1-(1,1-dimethylethyl)-3...	162	C12H18	000098-19-1	93
3		4-t-Butyl-o-xylene	162	C12H18	007397-06-0	93
4		Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	91
5		Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

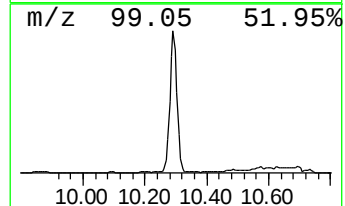
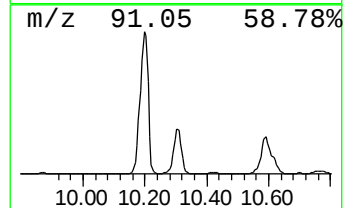
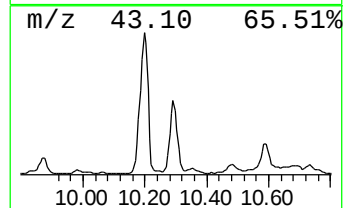
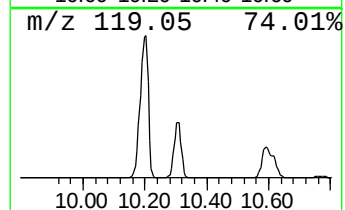
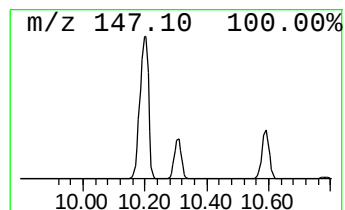
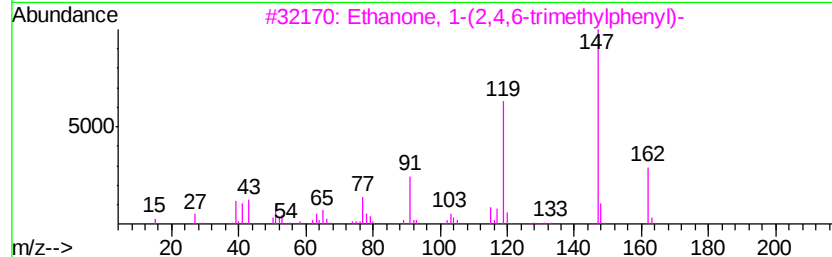
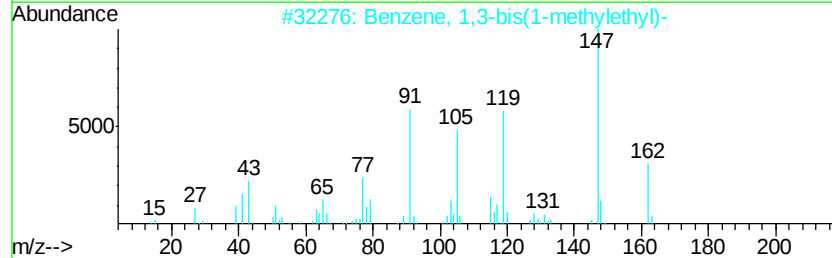
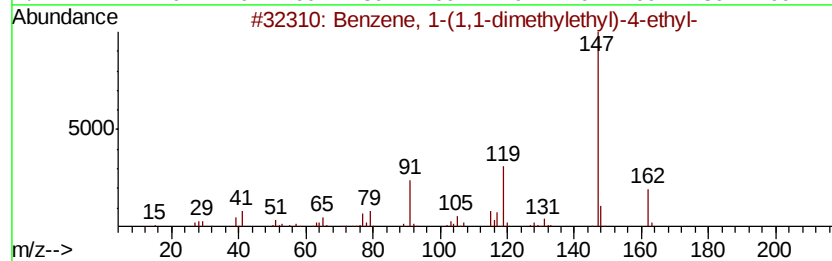
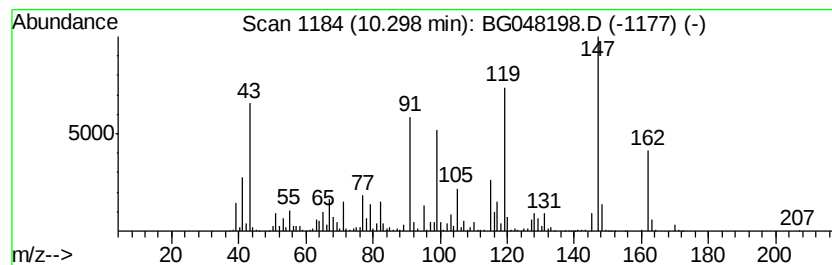
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Benzene, 1-(1,1-dimethyleth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	144.24 ng/ul	4973890	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-(1,1-dimethylethyl)-4...	162	C12H18	007364-19-4	91
2		Benzene, 1,3-bis(1-methylethyl)-	162	C12H18	000099-62-7	83
3		Ethanone, 1-(2,4,6-trimethylphen...	162	C11H14O	001667-01-2	81
4		4-t-Butyl-o-xylene	162	C12H18	007397-06-0	81
5		Ethanone, 1-(2,4,6-trimethylphen...	162	C11H14O	001667-01-2	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

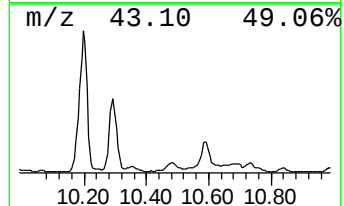
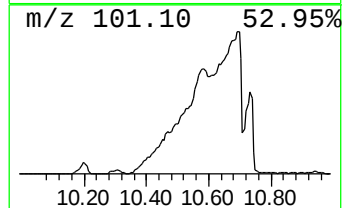
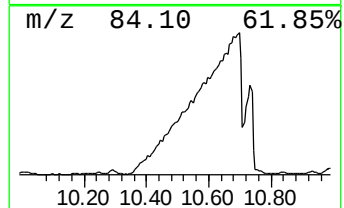
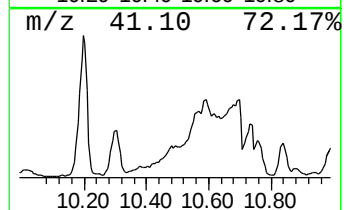
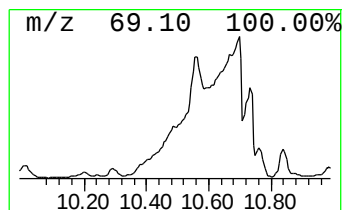
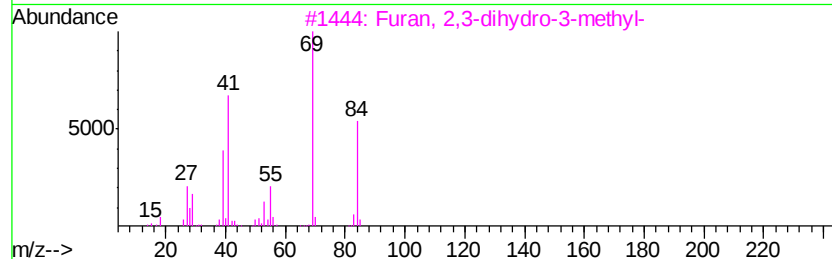
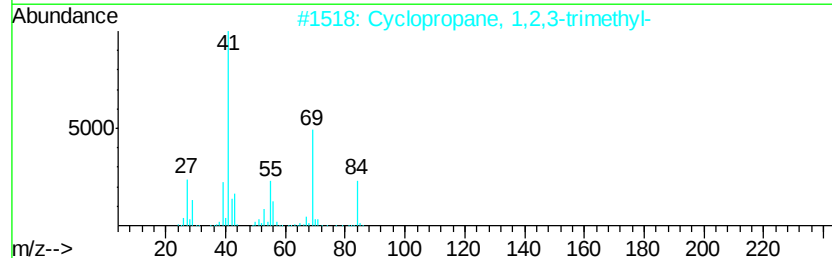
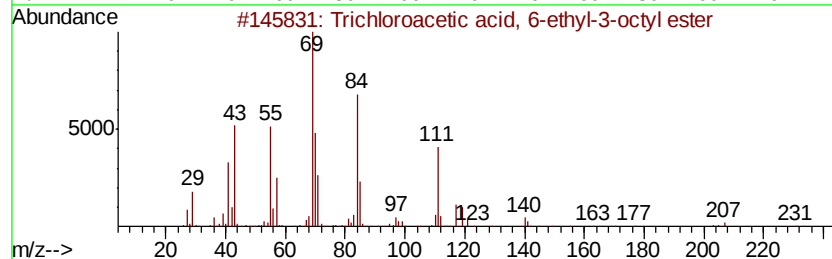
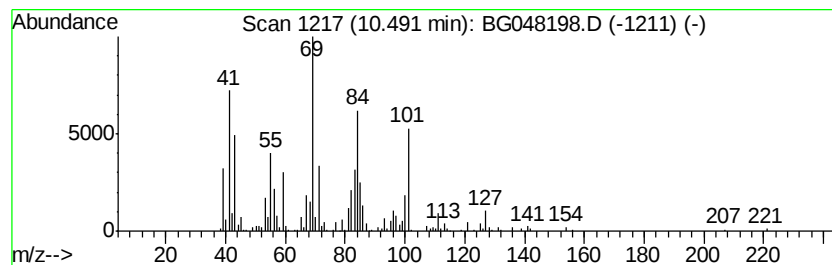
Instrument :
BNA_G
ClientSampleId :
DBGR3

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 unknown-08 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.49	14.43 ng/ul	497516	Naphthalene-d8	10.94		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trichloroacetic acid, 6-ethyl-3-...	302	C12H21Cl3O2	1000282-57-4	47
2		Cyclopropane, 1,2,3-trimethyl-	84	C6H12	042984-19-0	38
3		Furan, 2,3-dihydro-3-methyl-	84	C5H8O	001708-27-6	38
4		2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	38
5		2-Pentene, 3-methyl-	84	C6H12	000922-61-2	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
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Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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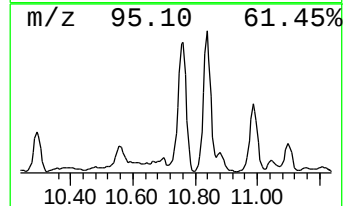
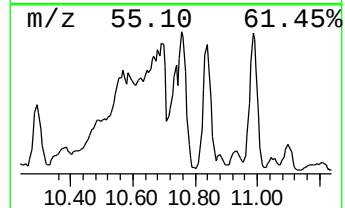
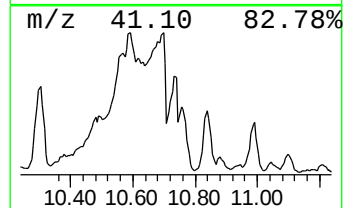
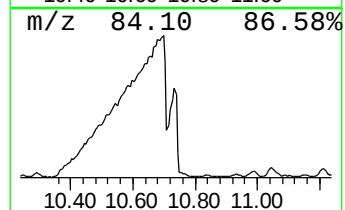
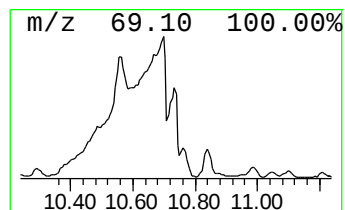
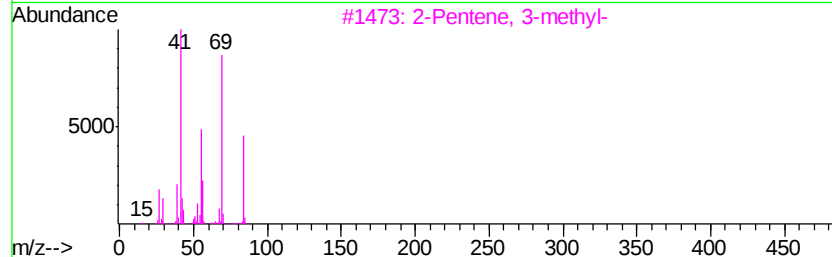
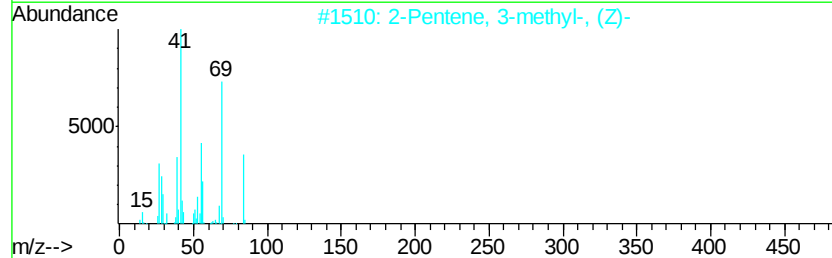
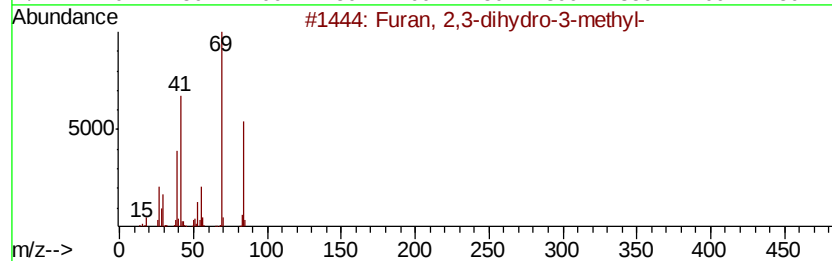
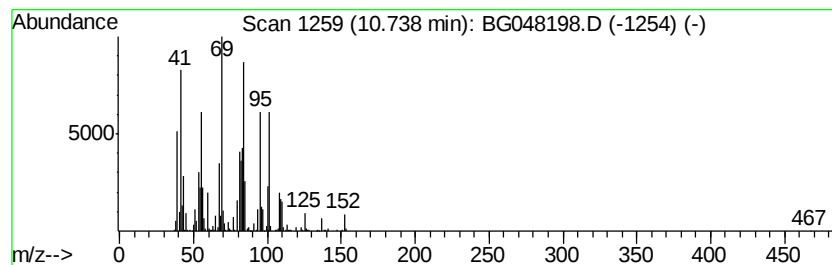
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 unknown-09 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.74	20.22 ng/ul	697158	Napthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Furan, 2,3-dihydro-3-methyl-	84	C5H8O	001708-27-6	35
2		2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	35
3		2-Pentene, 3-methyl-	84	C6H12	000922-61-2	35
4		3-Hexene, (E)-	84	C6H12	013269-52-8	35
5		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
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ALS Vial : 9 Sample Multiplier: 1

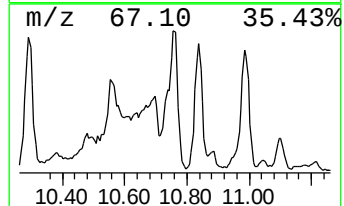
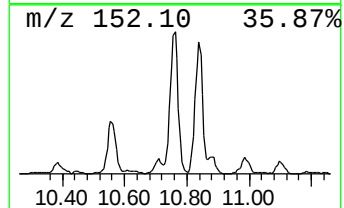
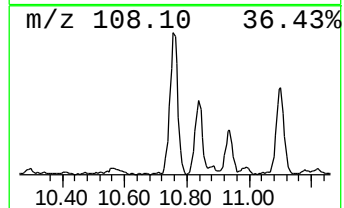
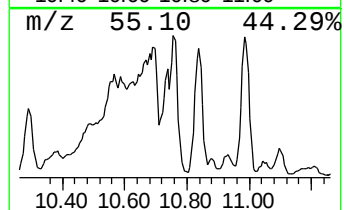
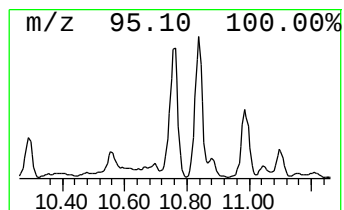
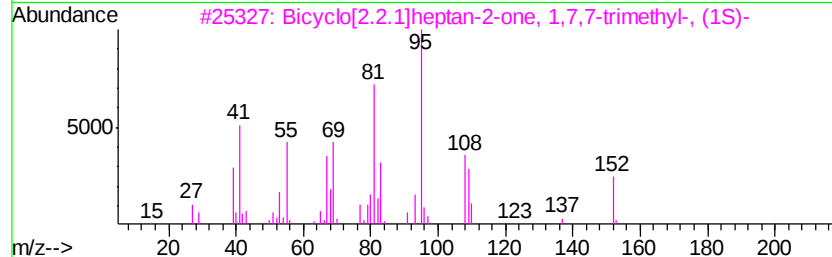
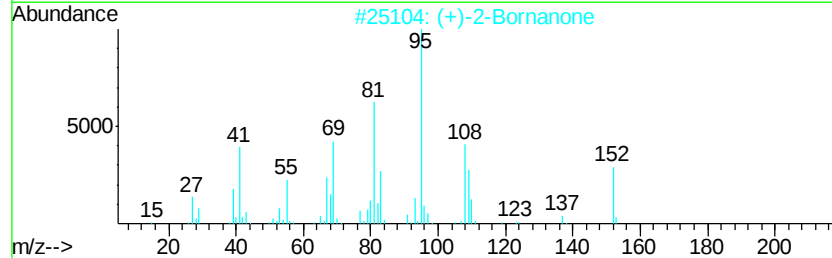
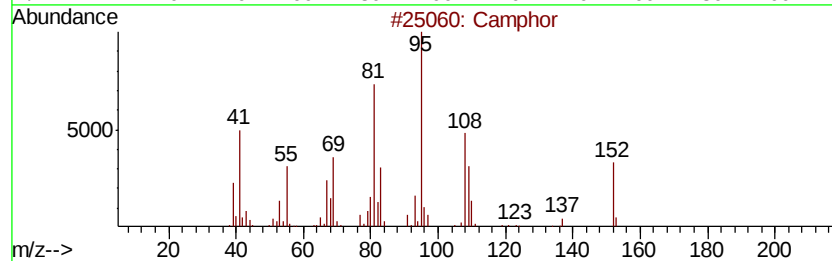
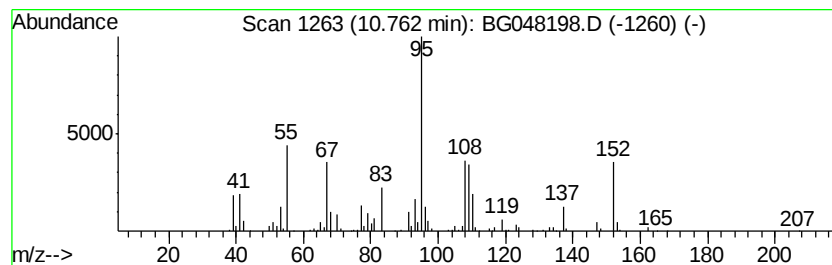
Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 Camphor Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	44.12 ng/ul	1521280	Napthalene-d8	10.94
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Camphor		152 C10H16O	000076-22-2 87
2	(+)-2-Bornanone		152 C10H16O	000464-49-3 86
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...		152 C10H16O	000464-48-2 83
4	Camphor		152 C10H16O	000076-22-2 83
5	(+)-2-Bornanone		152 C10H16O	000464-49-3 74



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Data File : BG048198.D
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Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

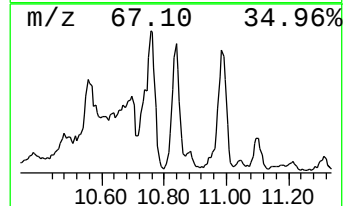
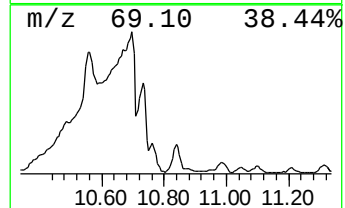
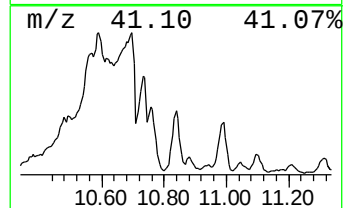
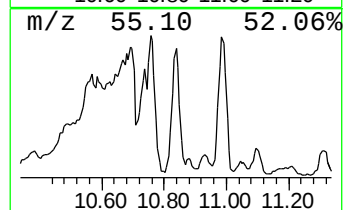
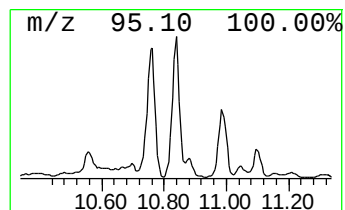
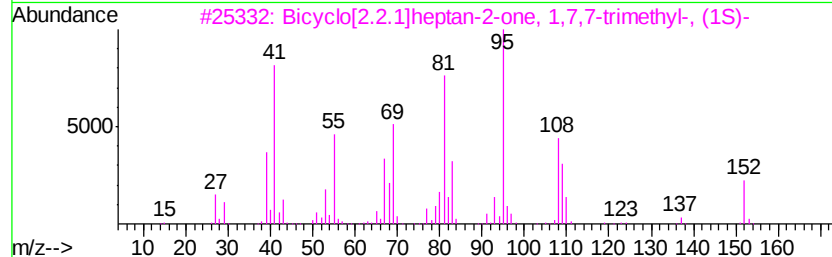
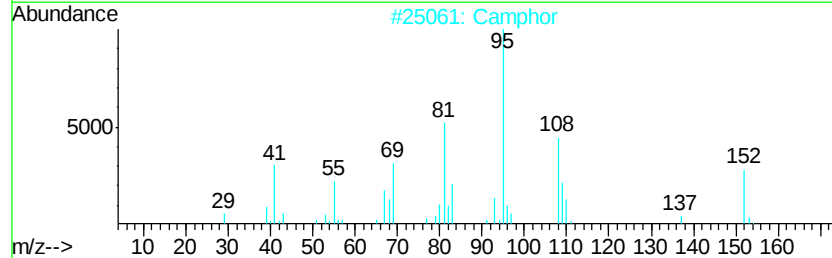
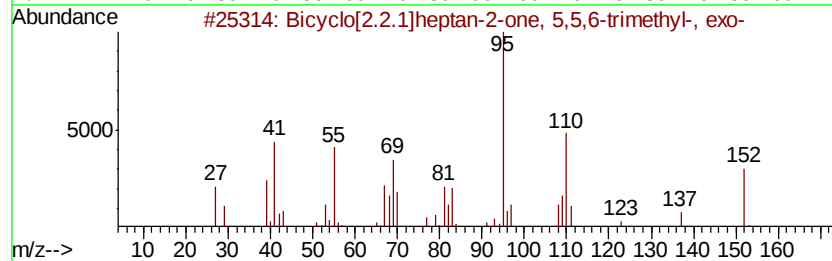
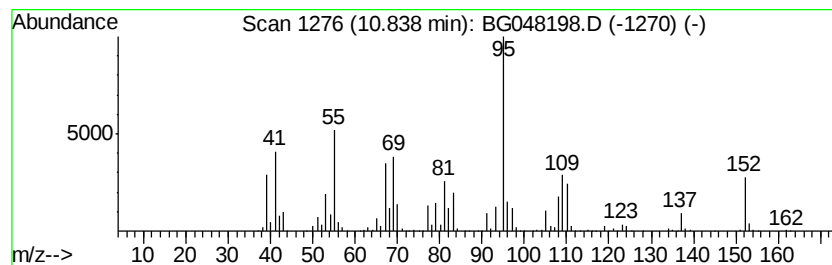
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	45.51 ng/ul	1569470	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptan-2-one, 5,5,...	152	C10H16O	003649-86-3	80
2		Camphor	152	C10H16O	000076-22-2	74
3		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	64
4		(+)-2-Bornanone	152	C10H16O	000464-49-3	64
5		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
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Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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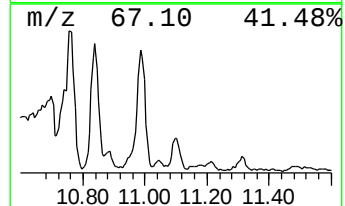
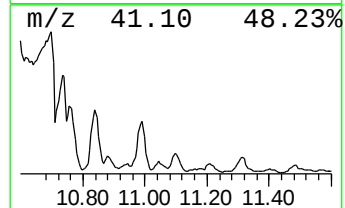
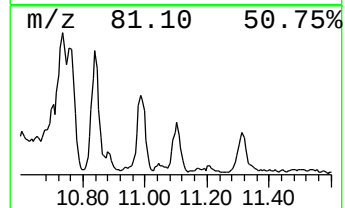
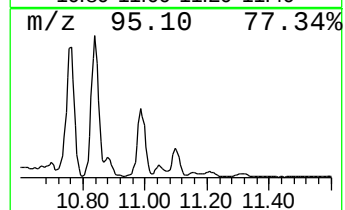
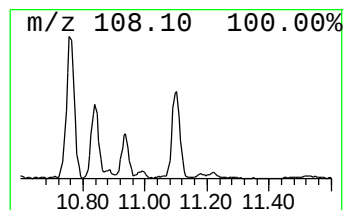
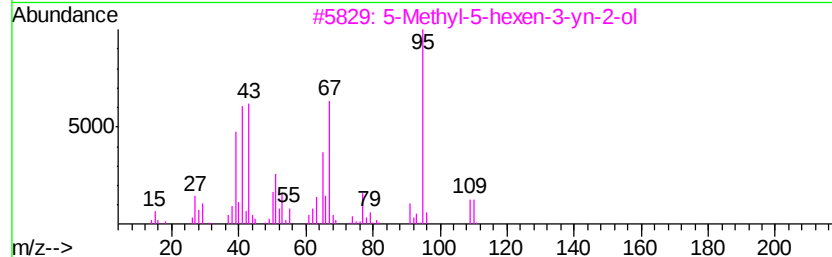
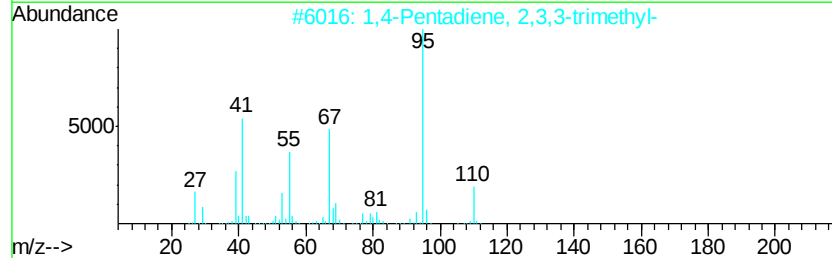
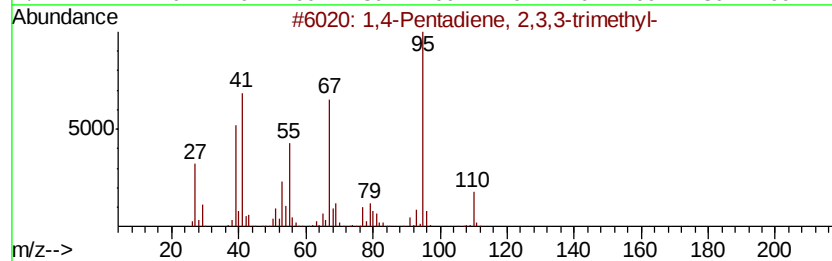
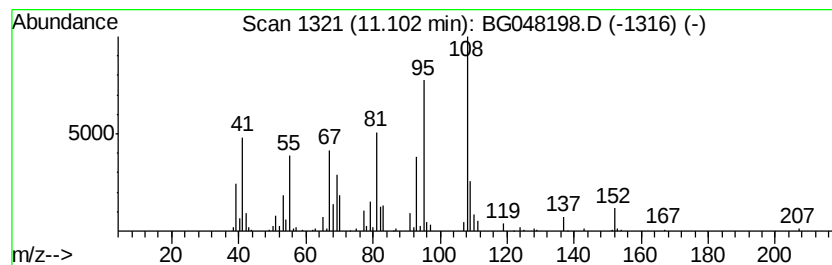
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 unknown-10 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	11.05 ng/ul	381017	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	30
2		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	30
3		5-Methyl-5-hexen-3-yn-2-ol	110	C7H10O	068017-33-4	30
4		Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	30
5		2,4-Hexadiene, 2,3-dimethyl-	110	C8H14	005678-98-8	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
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ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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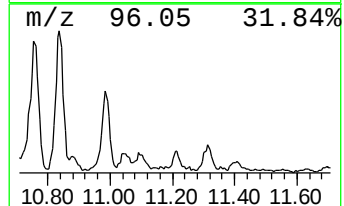
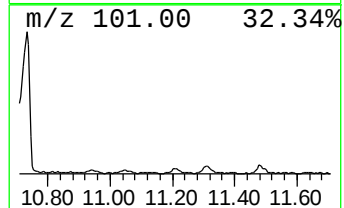
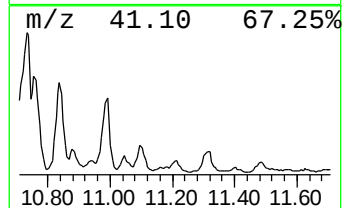
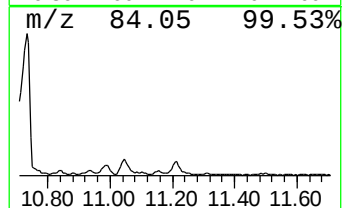
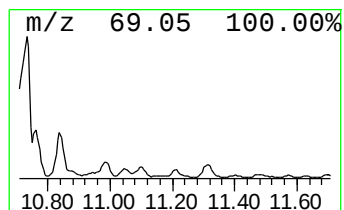
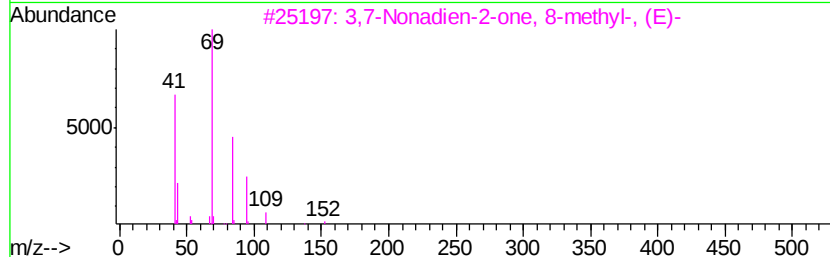
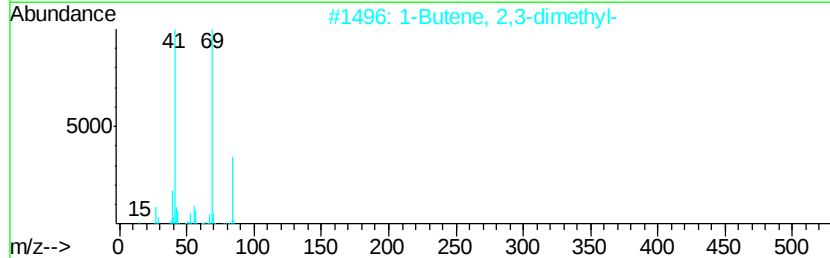
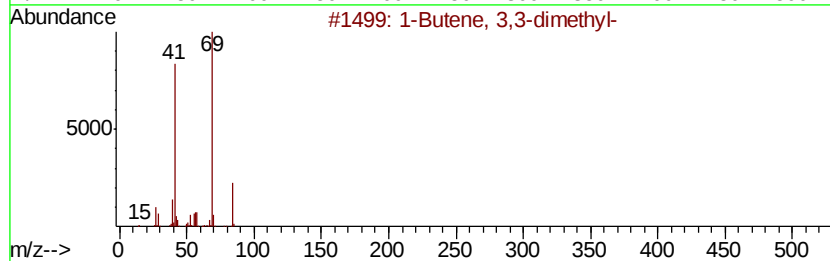
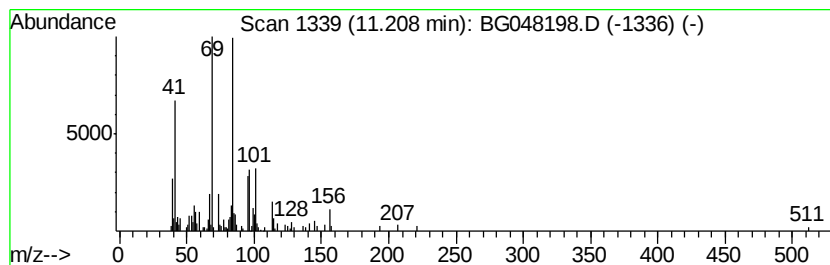
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.21	4.55 ng/ul	156878	Napthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butene, 3,3-dimethyl-	84	C6H12	000558-37-2	49
2			1-Butene, 2,3-dimethyl-	84	C6H12	000563-78-0	49
3			3,7-Nonadien-2-one, 8-methyl-, (E)-	152	C10H16O	035408-14-1	35
4			2,6-Dimethyl-2-trans-6-octadiene	138	C10H18	002609-23-6	27
5			2-Isopropenyl-5-methylhex-4-enal	152	C10H16O	075697-98-2	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

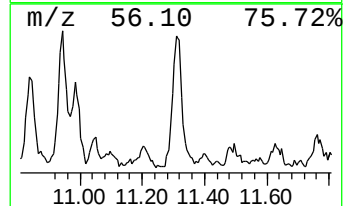
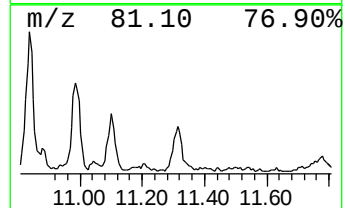
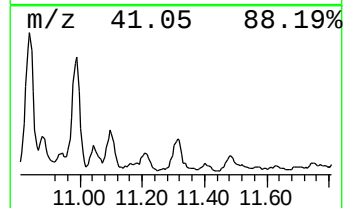
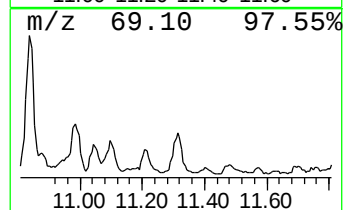
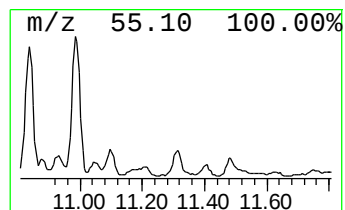
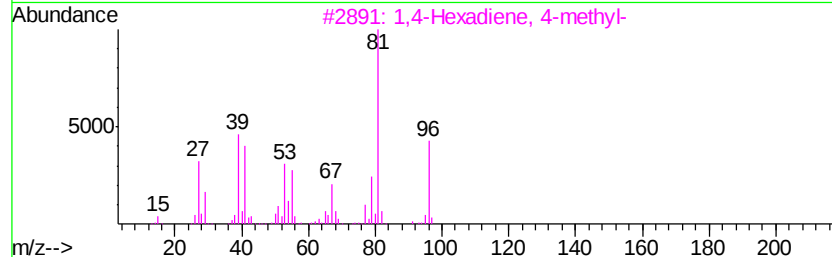
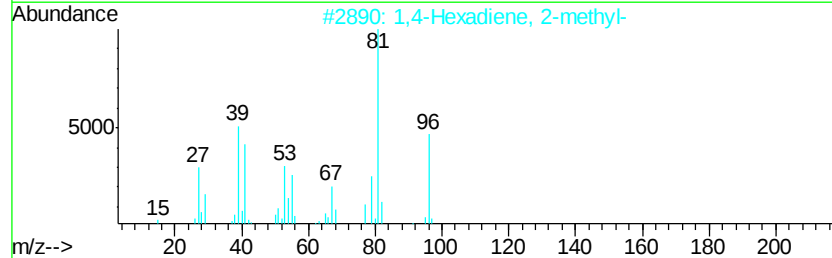
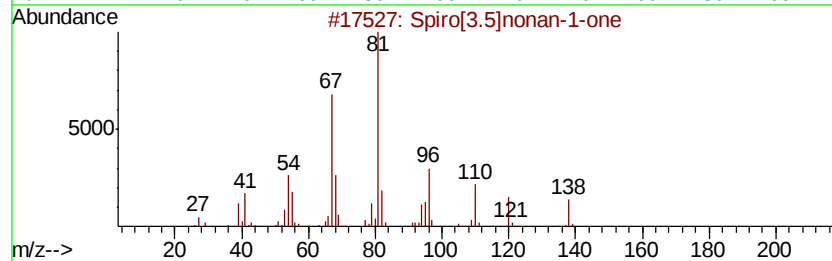
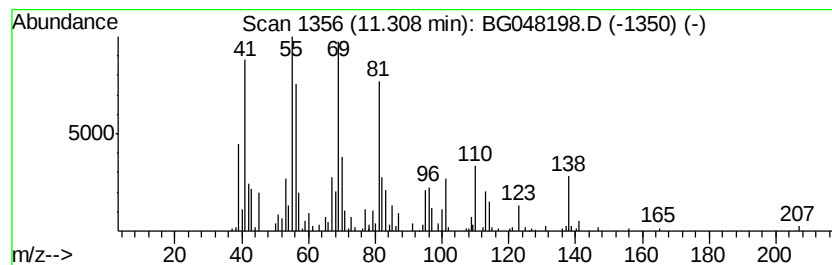
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 unknown-11 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.31	9.70 ng/ul	334582	Naphthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Spiro[3.5]nonan-1-one	138	C9H14O	029800-45-1	41
2			1,4-Hexadiene, 2-methyl-	96	C7H12	001119-14-8	35
3			1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	35
4			Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	35
5			2,7-Dioxatricyclo[4.3.1.0(3,8)]d...	138	C8H10O2	075568-65-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

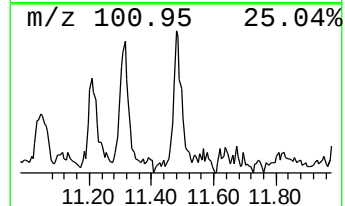
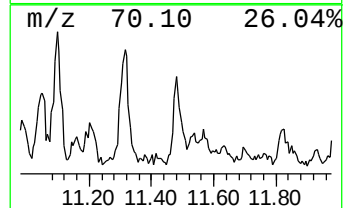
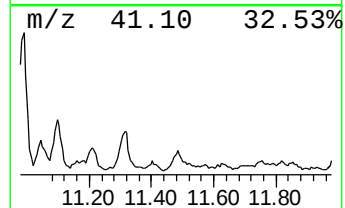
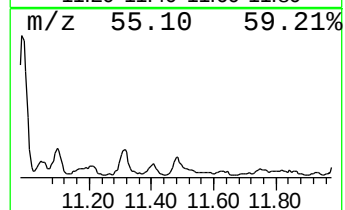
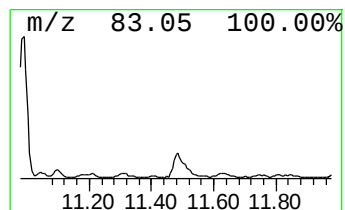
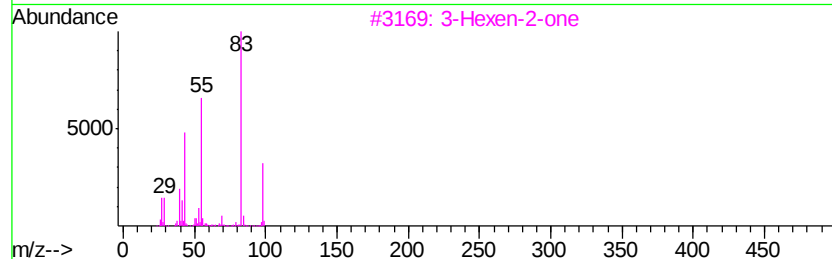
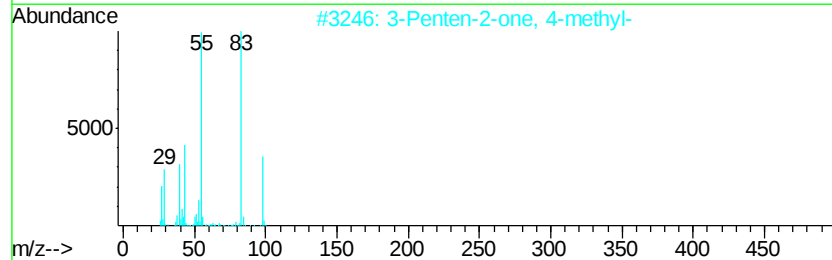
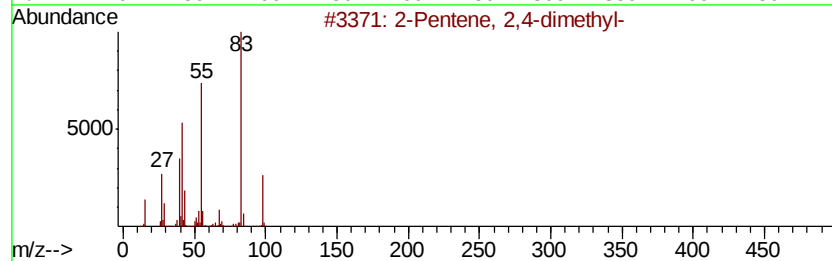
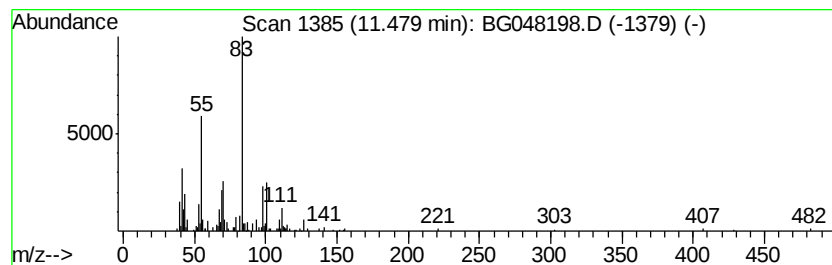
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	7.99 ng/ul	275568	Napthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	53
2		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	49
3		3-Hexen-2-one	98	C6H10O	000763-93-9	49
4		3-Hexen-2-one	98	C6H10O	000763-93-9	49
5		1H-Pyrazole, 4,5-dihydro-4,5-dim...	98	C5H10N2	028019-94-5	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

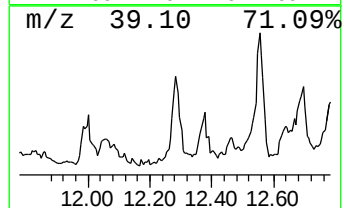
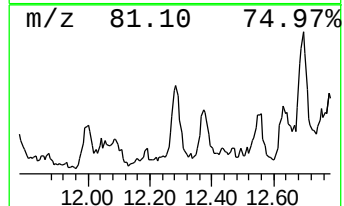
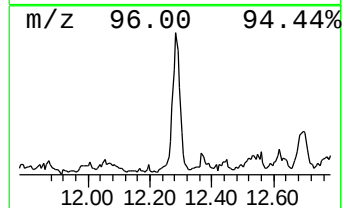
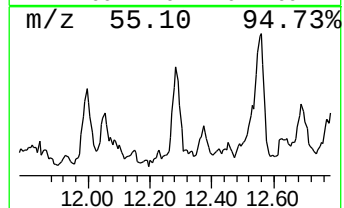
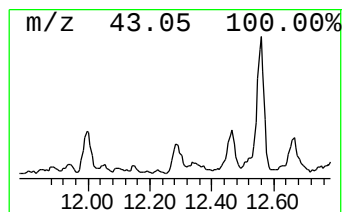
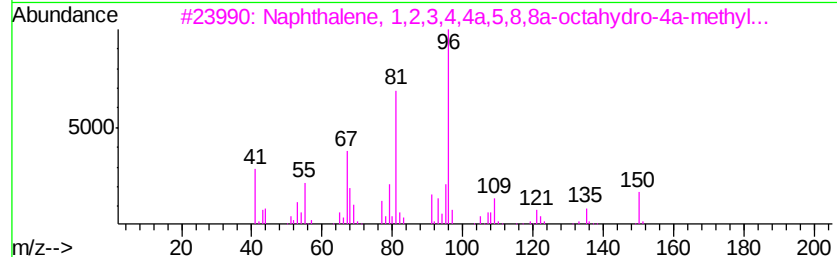
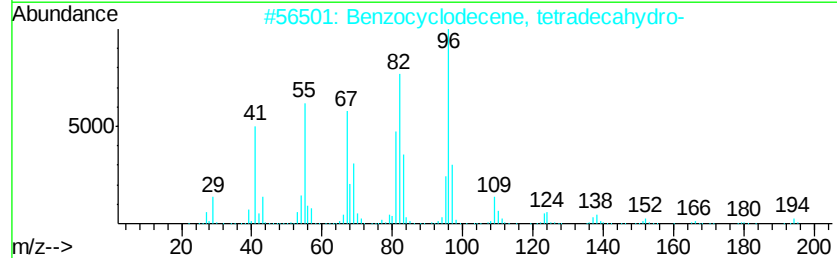
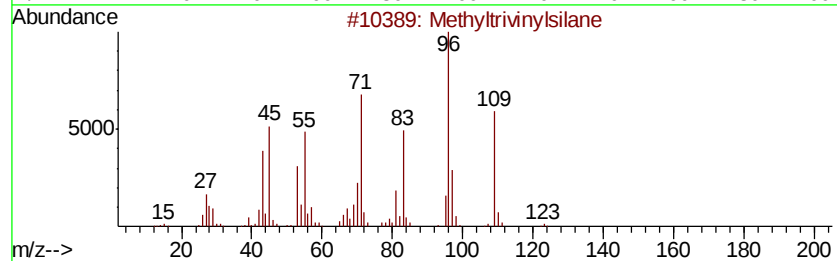
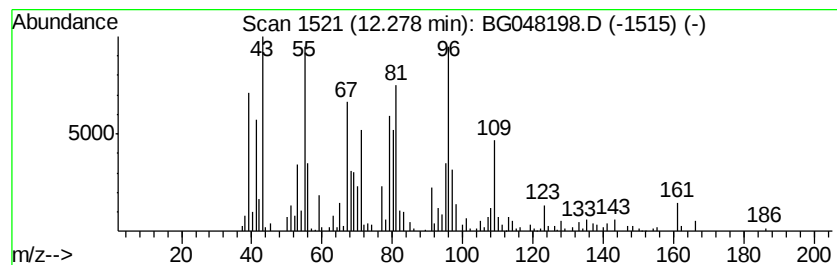
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 unknown-12 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	11.16 ng/ul	384935	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyltrivinylsilane	124	C7H12Si	018244-95-6	43
2		Benzocyclodecene, tetradecahydro-	194	C14H26	061142-06-1	38
3		Naphthalene, 1,2,3,4,4a,5,8,8a-o...	150	C11H18	021789-56-0	38
4		(4-Methyl-pent-3-enyl)-cyclohexane	166	C12H22	1000185-19-1	35
5		Cyclohexane, 1-ethenyl-2-methyl-...	124	C9H16	034780-45-5	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

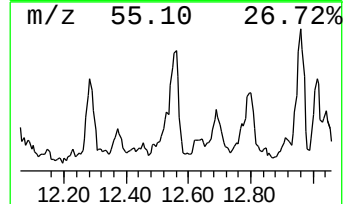
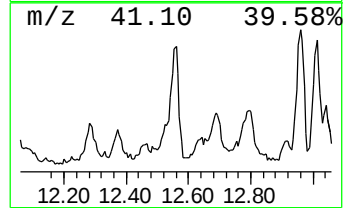
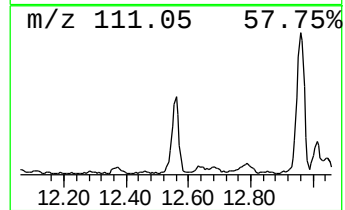
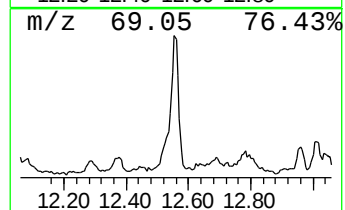
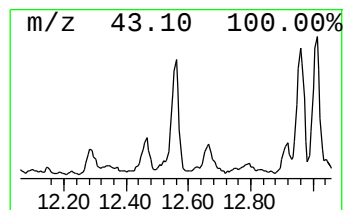
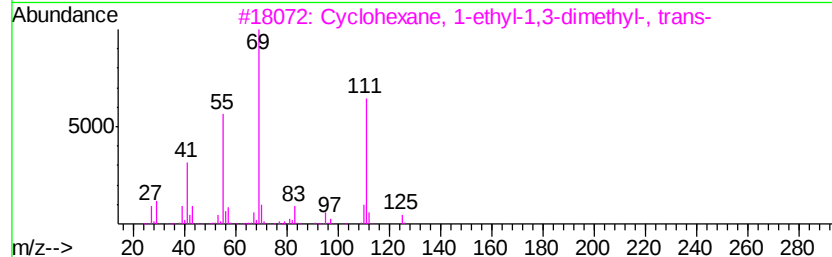
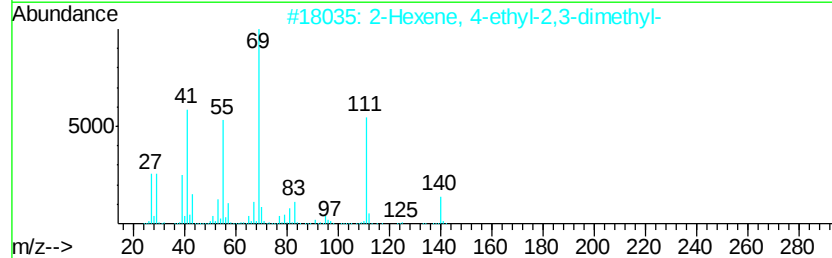
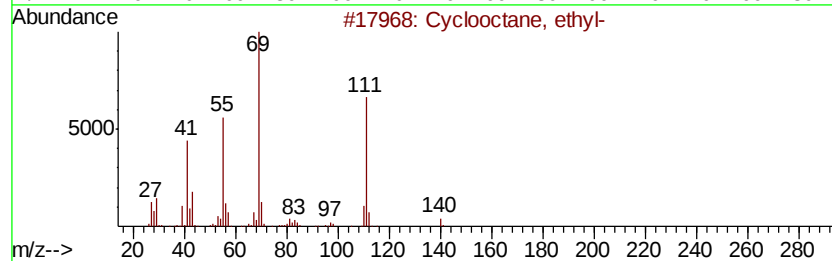
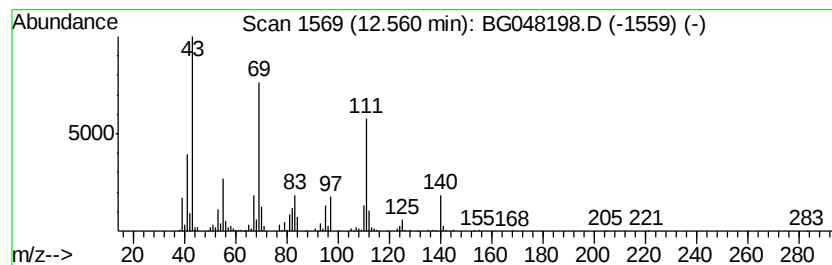
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 (DEL) Alkane: Cyclic12.56 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	20.33 ng/ul	700984	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclooctane, ethyl-	140	C10H20	013152-02-8	64
2		2-Hexene, 4-ethyl-2,3-dimethyl-	140	C10H20	1000149-19-7	64
3		Cyclohexane, 1-ethyl-1,3-dimethyl...	140	C10H20	062238-29-3	64
4		Cyclohexane, 1-ethyl-1,4-dimethyl...	140	C10H20	062238-30-6	64
5		Cyclohexane, 1-ethyl-1,4-dimethyl...	140	C10H20	062238-32-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

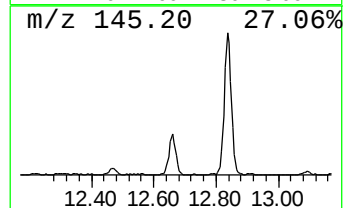
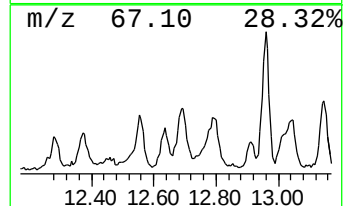
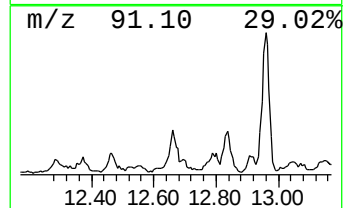
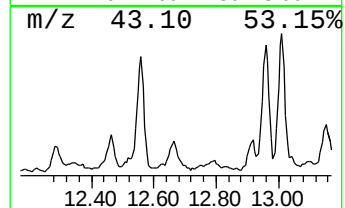
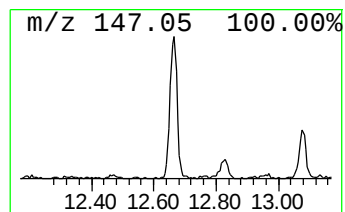
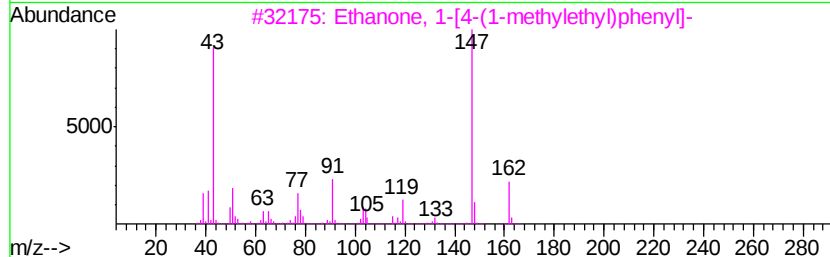
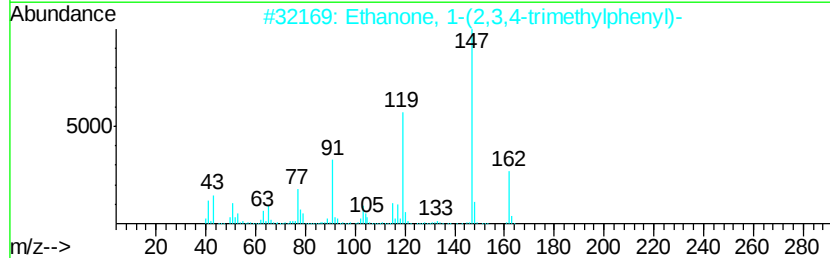
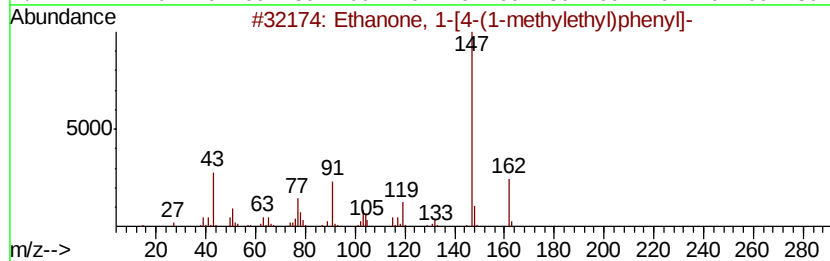
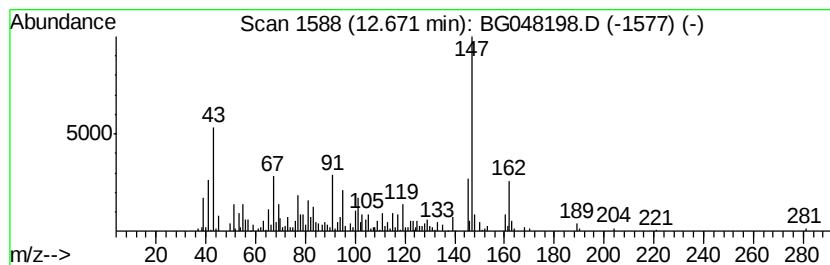
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 Ethanone, 1-[4-(1-methyleth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	11.08 ng/ul	382253	Naphthalene-d8	10.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-[4-(1-methylethyl)ph...	162	C11H14O	000645-13-6	70
2			Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	64
3			Ethanone, 1-[4-(1-methylethyl)ph...	162	C11H14O	000645-13-6	62
4			Benzene, hexamethyl-	162	C12H18	000087-85-4	58
5			5-Hydroxy-3-methyl-1-indanone	162	C10H10O2	057878-30-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

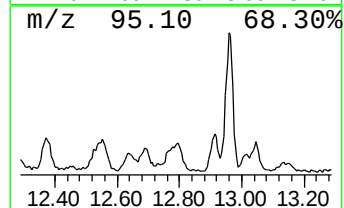
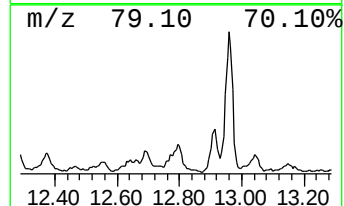
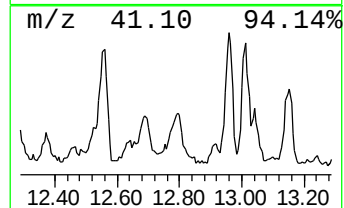
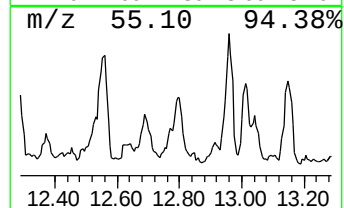
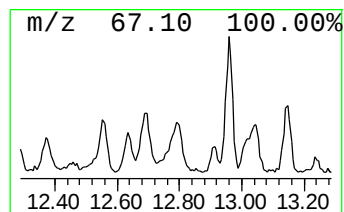
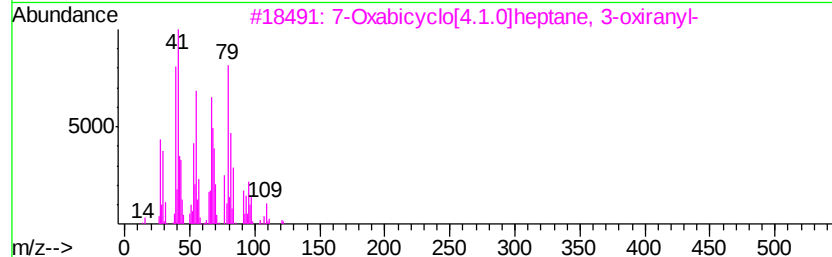
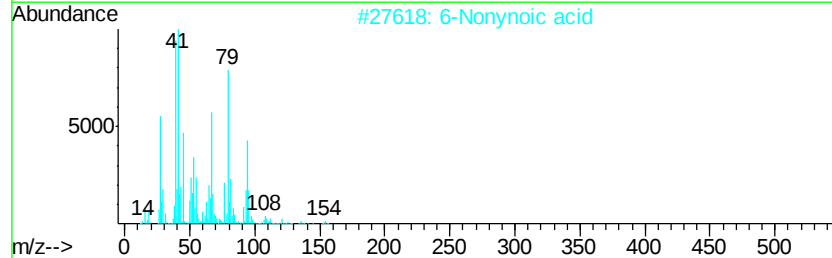
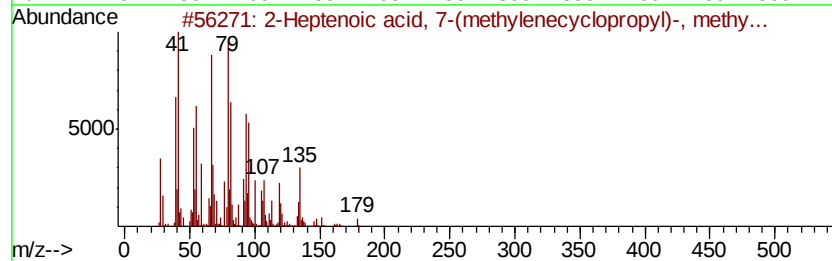
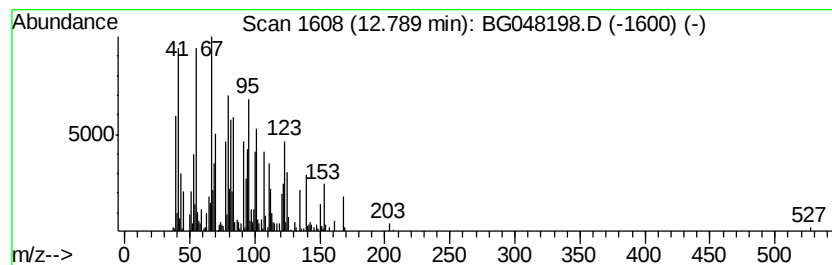
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 unknown-13 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	11.52 ng/ul	397386	Naphthalene-d8	10.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Heptenoic acid, 7-(methylenecyclopropyl)-, methyl ester	194	C12H18O2	1000158-04-1	43
2		6-Nonynoic acid	154	C9H14O2	056630-31-0	38
3		7-Oxabicyclo[4.1.0]heptane, 3-oxiranyl-	140	C8H12O2	000106-87-6	38
4		6-Methyl-2-heptyne	110	C8H14	051065-64-6	35
5		2-Hexyn-1-ol	98	C6H10O	000764-60-3	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

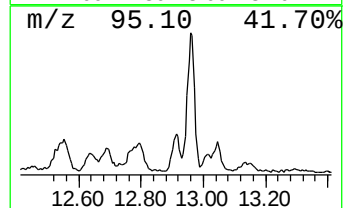
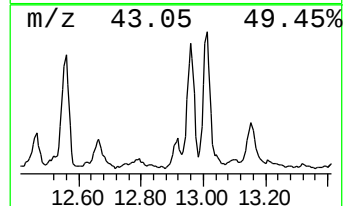
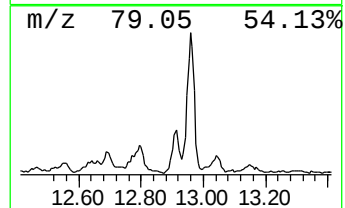
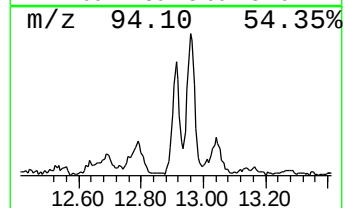
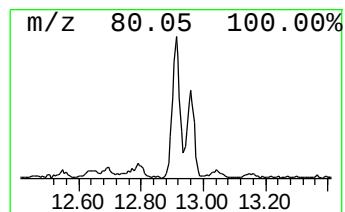
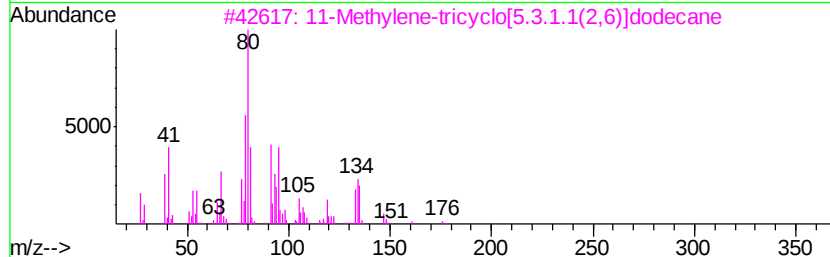
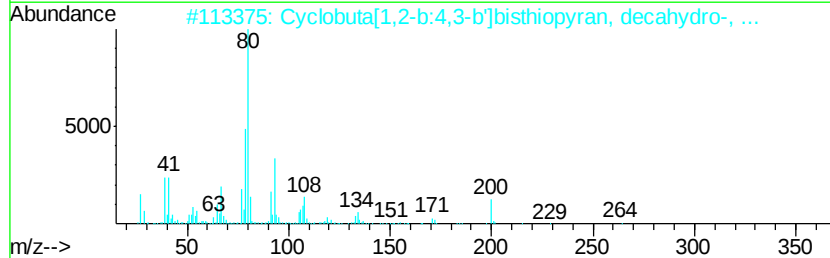
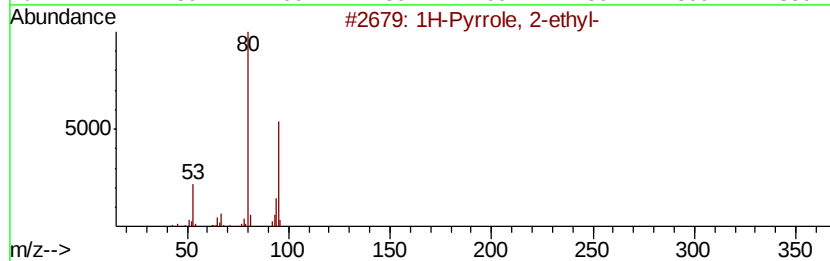
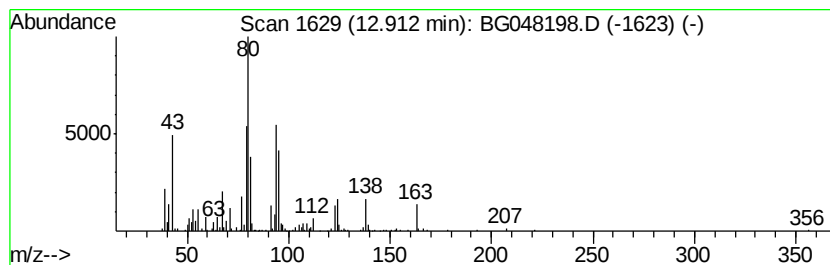
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 1H-Pyrrole, 2-ethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	10.23 ng/ul	301533	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrrole, 2-ethyl-	95	C6H9N	001551-06-0	64
2		Cyclobuta[1,2-b:4,3-b']bisthiopy...	264	C10H16O4S2	074630-21-0	59
3		11-Methylene-tricyclo[5.3.1.1(2,...	176	C13H20	1000193-95-2	56
4		Bicyclo[2.2.1]hept-2-ene, 2-methyl-	108	C8H12	000694-92-8	53
5		Bicyclo[3.2.1]octan-2-one	124	C8H12O	005019-82-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBG3

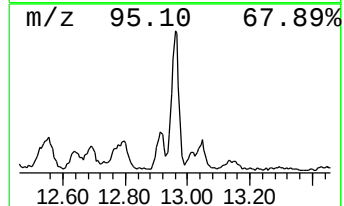
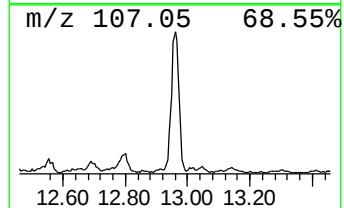
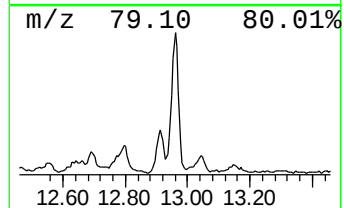
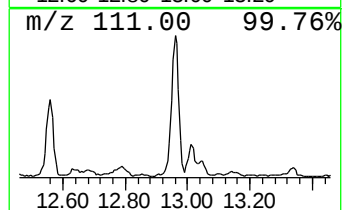
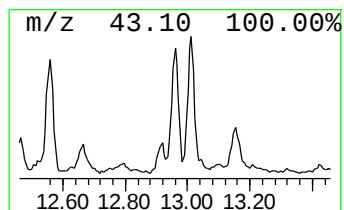
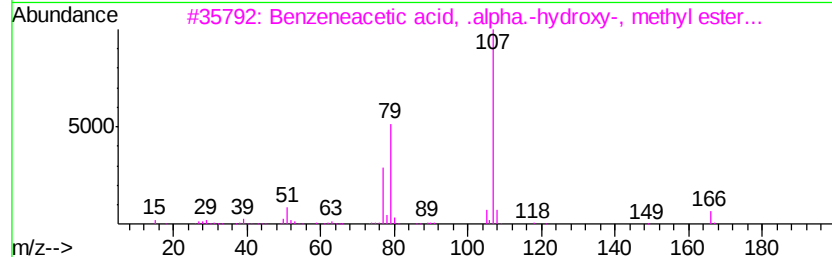
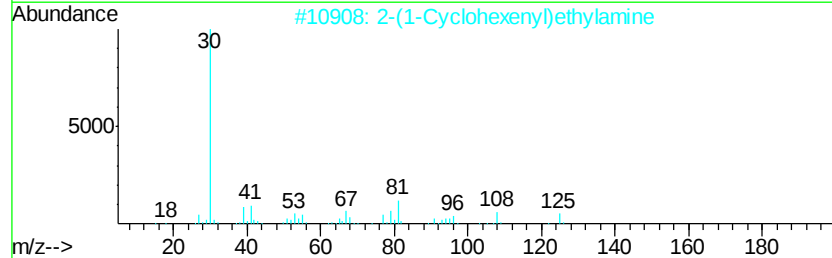
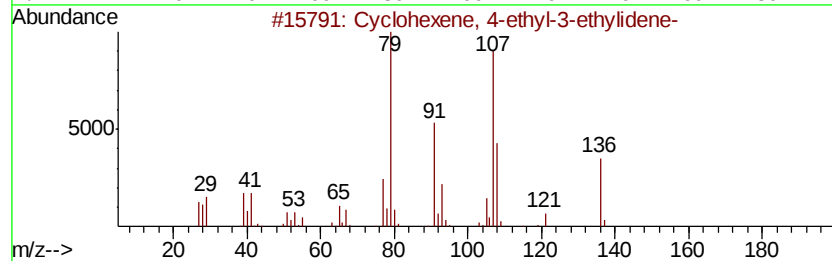
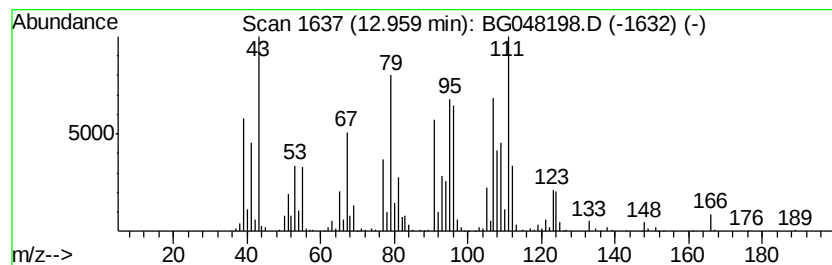
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 unknown-14 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	41.13 ng/ul	1212840	Acenaphthene-d10	14.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 4-ethyl-3-ethylidene-	136	C10H16	061233-77-0	22
2			2-(1-Cyclohexenyl)ethylamine	125	C8H15N	003399-73-3	22
3			Benzeneacetic acid, .alpha.-hvd...	166	C9H10O3	021210-43-5	18
4			1,3-trans,5-cis-Octatriene	108	C8H12	040087-61-4	18
5			E,Z-4-Ethylidenecyclohexene	108	C8H12	016631-66-6	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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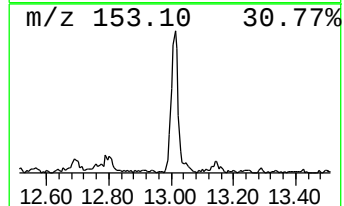
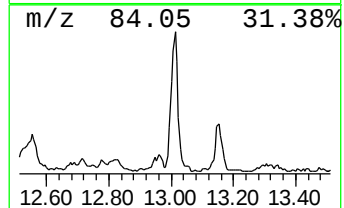
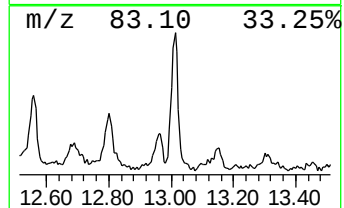
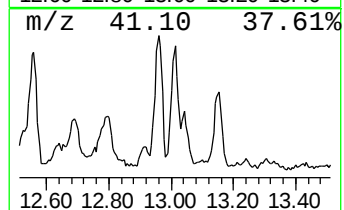
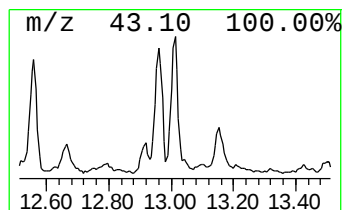
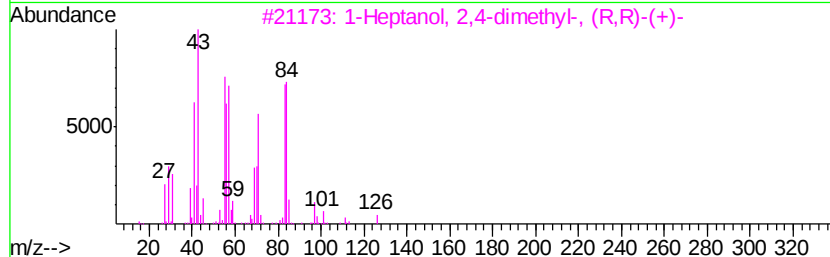
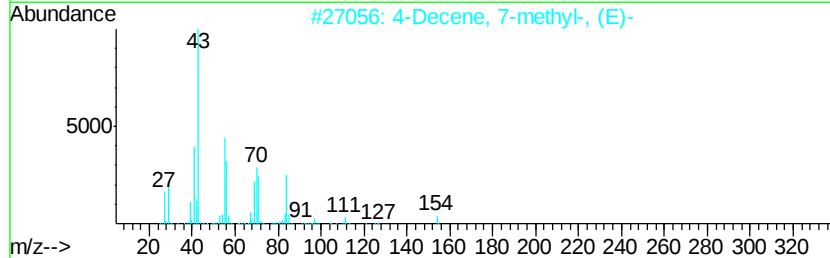
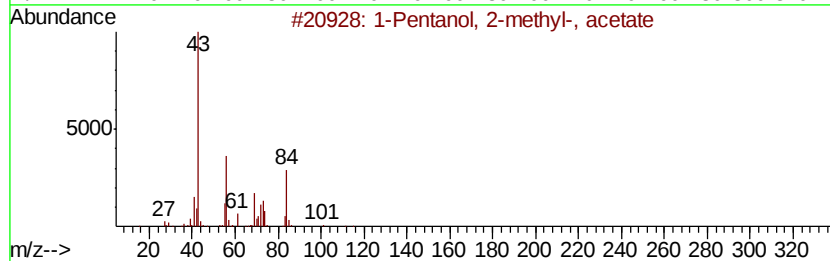
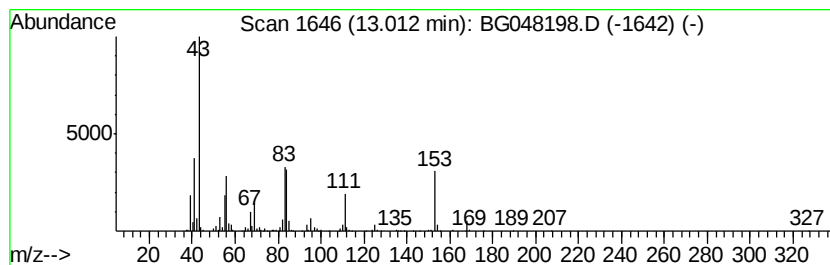
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 unknown-15 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	14.16 ng/ul	417631	Acenaphthene-d10	14.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentanol, 2-methyl-, acetate	144	C8H16O2	007789-99-3	25
2			4-Decene, 7-methyl-, (E)-	154	C11H22	062338-48-1	25
3			1-Heptanol, 2,4-dimethyl-, (R,R)-...	144	C9H20O	018450-73-2	23
4			1-[1,2,4]Triazol-1-yl-ethanone	111	C4H5N3O	015625-88-4	14
5			dl-Lysine	146	C6H14N2O2	000070-54-2	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
 Data File : BG048198.D
 Acq On : 18 Feb 2021 16:44
 Operator : CG/JU
 Sample : M1429-03
 Misc :
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Instrument :
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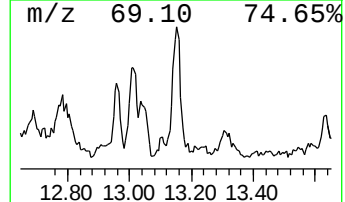
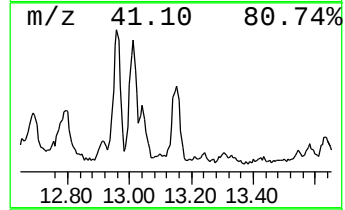
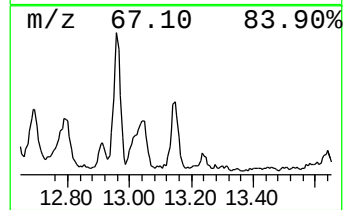
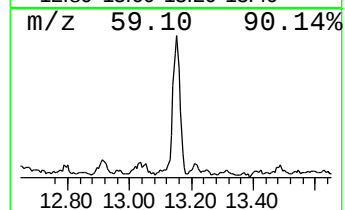
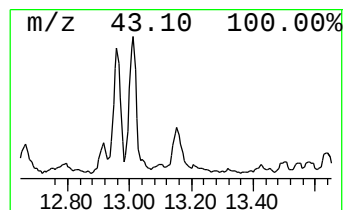
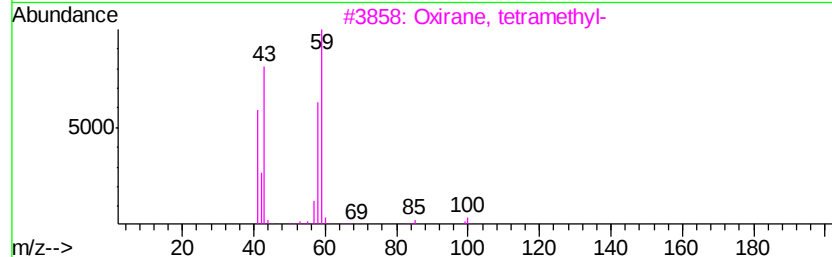
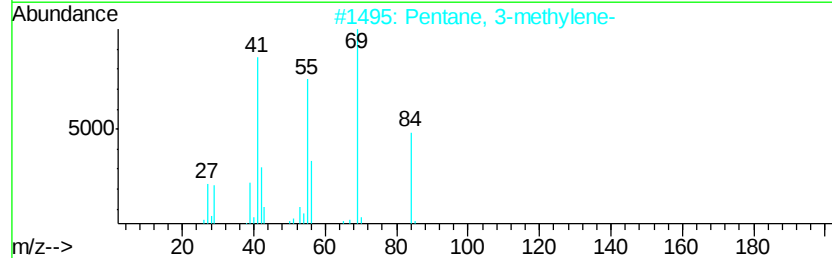
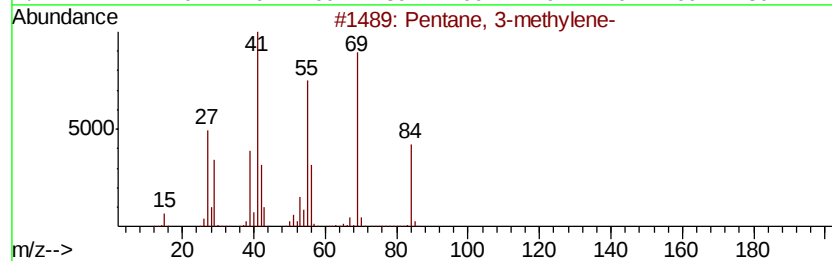
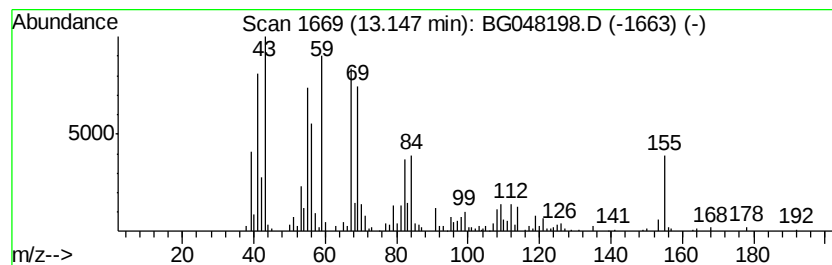
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 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	13.99 ng/ul	412464	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methylene-	84	C6H12	000760-21-4	25
2		Pentane, 3-methylene-	84	C6H12	000760-21-4	25
3		Oxirane, tetramethyl-	100	C6H12O	005076-20-0	25
4		Guanidine	59	CH5N3	000113-00-8	22
5		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
Acq On : 18 Feb 2021 16:44
Operator : CG/JU
Sample : M1429-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBGR3

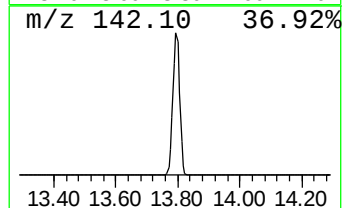
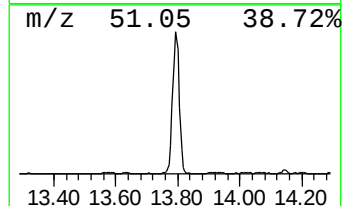
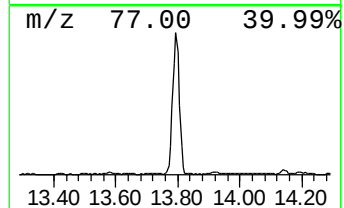
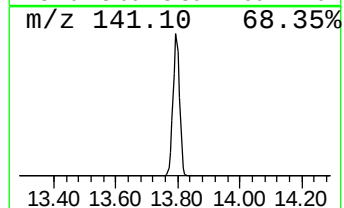
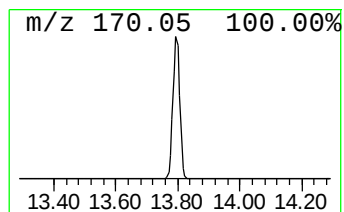
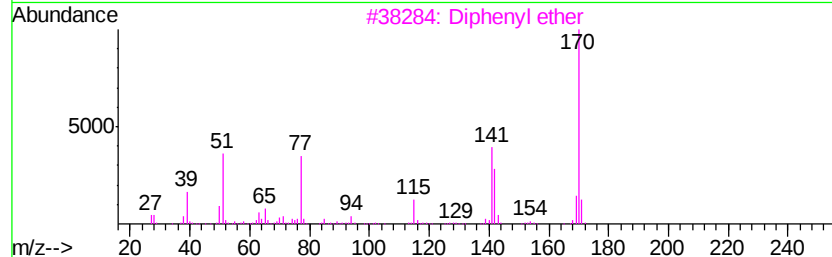
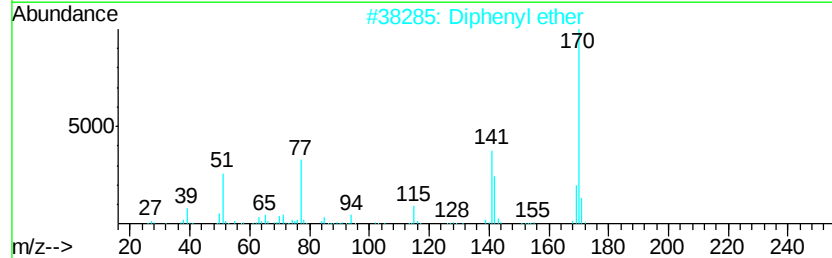
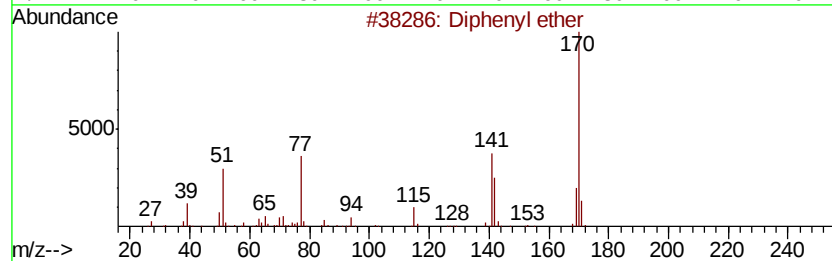
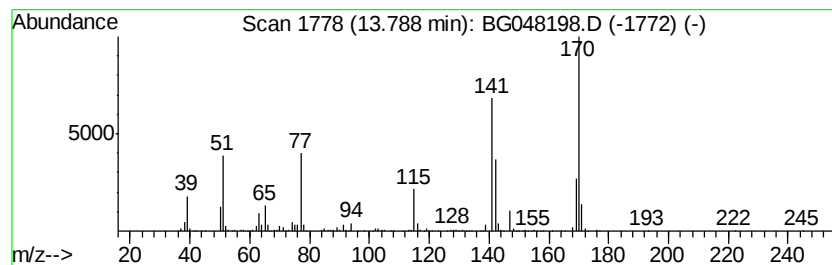
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 Diphenyl ether Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	109.86 ng/ul	3239720	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diphenyl ether	170	C12H10O	000101-84-8	87
2		Diphenyl ether	170	C12H10O	000101-84-8	81
3		Diphenyl ether	170	C12H10O	000101-84-8	81
4		Diphenyl ether	170	C12H10O	000101-84-8	64
5		1H,3H-Naphtho[1,8-cd]pyran	170	C12H10O	000203-84-9	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021821\
Data File : BG048198.D
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Sample : M1429-03
Misc :
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Instrument :
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ClientSampleId :
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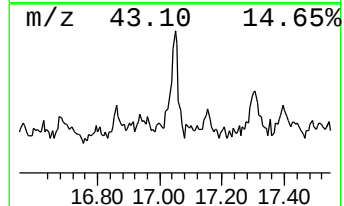
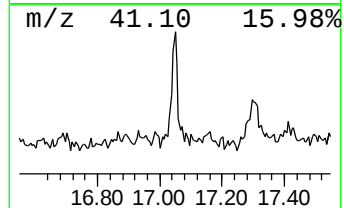
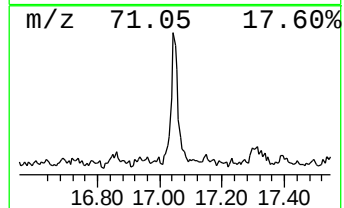
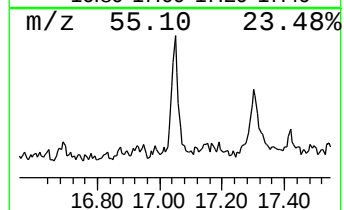
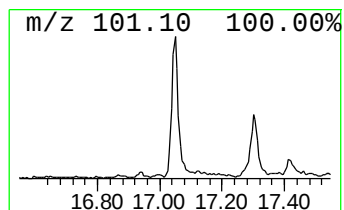
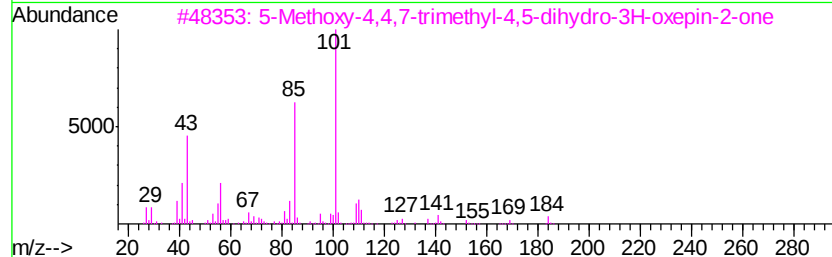
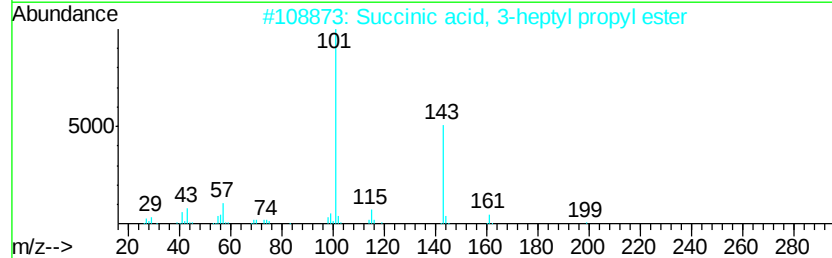
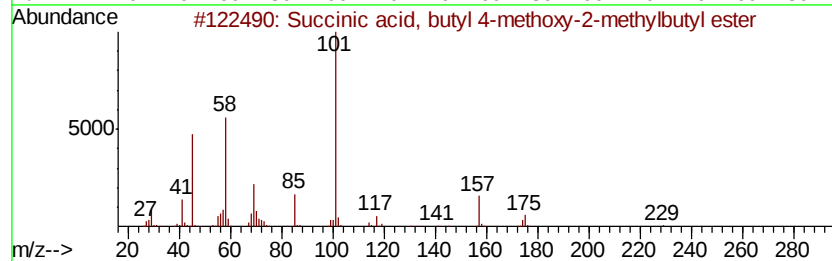
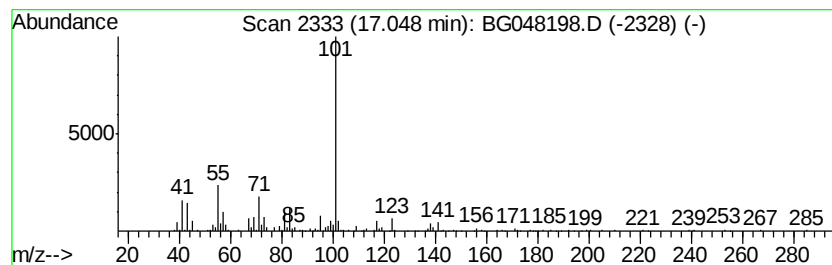
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 Succinic acid, butyl 4-meth... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.05	7.70 ng/ul	264117	Phenanthrene-d10	17.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Succinic acid, butyl 4-methoxy-2...	274	C14H26O5	1000330-23-2	64
2		Succinic acid, 3-heptyl propyl e...	258	C14H26O4	1000349-48-4	59
3		5-Methoxy-4,4,7-trimethyl-4,5-di...	184	C10H16O3	1000187-23-9	56
4		1,3-Dioxepane, 2-pentadecyl-	312	C20H40O2	041563-29-5	53
5		Succinic acid, hexadecyl 4-metho...	442	C26H50O5	1000330-24-2	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021821\
 Data File : BG048198.D
 Acq On : 18 Feb 2021 16:44
 Operator : CG/JU
 Sample : M1429-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexanol, 1-m...	6.20	76.7	ng/ul	1194380	1	8.11	311400	20.0
2-Pentanol, 3-eth...	6.44	3.4	ng/ul	53488	1	8.11	311400	20.0
Benzene, (1-methy...	6.59	16.7	ng/ul	260146	1	8.11	311400	20.0
unknown-01	7.55	10.2	ng/ul	159269	1	8.11	311400	20.0
unknown-02	7.71	6.7	ng/ul	103533	1	8.11	311400	20.0
(DEL) Alkane: Cyc...	7.75	9.0	ng/ul	139831	1	8.11	311400	20.0
unknown-03	7.94	5.0	ng/ul	77063	1	8.11	311400	20.0
unknown-04	9.14	10.4	ng/ul	162076	1	8.11	311400	20.0
unknown-05	9.38	65.8	ng/ul	1024510	1	8.11	311400	20.0
Benzene, 1-ethyl-...	9.45	28.0	ng/ul	435352	1	8.11	311400	20.0
unknown-06	9.63	6.5	ng/ul	222582	2	10.94	689683	20.0
Benzene, 1-ethyl-...	9.69	8.7	ng/ul	299344	2	10.94	689683	20.0
Buta-1,3-diene, 1...	9.79	8.3	ng/ul	284594	2	10.94	689683	20.0
(DEL) Alkane: Cyc...	9.83	6.4	ng/ul	219210	2	10.94	689683	20.0
unknown-07	9.87	19.6	ng/ul	675370	2	10.94	689683	20.0
Benzene, 1,3-bis(...	10.20	404.1	ng/ul	13934100	2	10.94	689683	20.0
Benzene, 1-(1,1-d...	10.30	144.2	ng/ul	4973890	2	10.94	689683	20.0
unknown-08	10.49	14.4	ng/ul	497516	2	10.94	689683	20.0
unknown-09	10.74	20.2	ng/ul	697158	2	10.94	689683	20.0
Camphor	10.76	44.1	ng/ul	1521280	2	10.94	689683	20.0
Bicyclo[2.2.1]hep...	10.84	45.5	ng/ul	1569470	2	10.94	689683	20.0
unknown-10	11.10	11.1	ng/ul	381017	2	10.94	689683	20.0
(DEL) Alkane: Str...	11.21	4.5	ng/ul	156878	2	10.94	689683	20.0
unknown-11	11.31	9.7	ng/ul	334582	2	10.94	689683	20.0
(DEL) Alkane: Str...	11.48	8.0	ng/ul	275568	2	10.94	689683	20.0
unknown-12	12.28	11.2	ng/ul	384935	2	10.94	689683	20.0
(DEL) Alkane: Cyc...	12.56	20.3	ng/ul	700984	2	10.94	689683	20.0
Ethanone, 1-[4-(1...	12.67	11.1	ng/ul	382253	2	10.94	689683	20.0
unknown-13	12.79	11.5	ng/ul	397386	2	10.94	689683	20.0
1H-Pyrrole, 2-ethyl-	12.91	10.2	ng/ul	301533	3	14.74	589773	20.0
unknown-14	12.96	41.1	ng/ul	1212840	3	14.74	589773	20.0
unknown-15	13.01	14.2	ng/ul	417631	3	14.74	589773	20.0
(DEL) Alkane: Str...	13.15	14.0	ng/ul	412464	3	14.74	589773	20.0
Diphenyl ether	13.79	109.9	ng/ul	3239720	3	14.74	589773	20.0
Succinic acid, bu...	17.05	7.7	ng/ul	264117	4	17.48	685865	20.0