

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
 Data File : BG049559.D
 Acq On : 29 Jul 2021 14:32
 Operator : CG/JU
 Sample : M3123-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0068

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049559.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.085	3	8	16	rBV2	88193	186903	45.86%	3.644%
2	3.156	16	20	35	rVB2	128264	222609	54.62%	4.340%
3	3.878	140	143	151	rVB4	5479	9479	2.33%	0.185%
4	4.889	313	315	318	rBV3	1210	1391	0.34%	0.027%
5	4.936	322	323	326	rBV	1751	1546	0.38%	0.030%
6	5.047	339	342	344	rBV2	1890	2051	0.50%	0.040%
7	5.212	368	370	376	rVB4	2418	3940	0.97%	0.077%
8	5.259	376	378	381	rBV2	1321	1790	0.44%	0.035%
9	5.594	432	435	436	rBV2	1325	1411	0.35%	0.028%
10	5.670	443	448	454	rVB4	2725	7304	1.79%	0.142%
11	6.040	507	511	515	rBV	5548	9322	2.29%	0.182%
12	6.234	540	544	545	rBV	1486	1416	0.35%	0.028%
13	6.669	616	618	621	rBV4	1478	1809	0.44%	0.035%
14	6.892	653	656	658	rBV	1673	1480	0.36%	0.029%
15	7.027	677	679	681	rBV2	2114	1630	0.40%	0.032%
16	7.068	685	686	689	rBV2	1596	1649	0.40%	0.032%
17	7.127	695	696	699	rVB2	3040	1630	0.40%	0.032%
18	7.203	702	709	716	rBV2	48772	95064	23.33%	1.853%
19	7.368	731	737	744	rBV2	28252	49851	12.23%	0.972%
20	7.579	766	773	779	rBV2	47662	83926	20.59%	1.636%
21	7.667	783	788	790	rBV3	4826	6447	1.58%	0.126%
22	8.049	844	853	863	rBV	86999	151467	37.17%	2.953%
23	8.689	959	962	965	rBV3	2120	2925	0.72%	0.057%
24	8.754	967	973	982	rBV	57757	99864	24.50%	1.947%
25	8.883	991	995	998	rBV5	3189	4284	1.05%	0.084%
26	9.012	1013	1017	1022	rBV	1086	2400	0.59%	0.047%
27	9.218	1044	1052	1058	rBV	49415	89118	21.87%	1.738%
28	9.271	1058	1061	1068	rVB4	8038	14569	3.57%	0.284%
29	9.365	1075	1077	1078	rBV4	1641	1574	0.39%	0.031%
30	9.618	1117	1120	1122	rBV4	1211	1703	0.42%	0.033%
31	9.694	1130	1133	1134	rBV3	1502	1653	0.41%	0.032%
32	9.782	1146	1148	1151	rVB2	1499	1503	0.37%	0.029%
33	9.829	1154	1156	1161	rVB	1440	1421	0.35%	0.028%
34	9.946	1169	1176	1186	rBV	46197	83314	20.44%	1.624%
35	10.187	1214	1217	1219	rBV	1216	1417	0.35%	0.028%
36	10.281	1231	1233	1238	rVB2	2094	2368	0.58%	0.046%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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37	10.387	1243	1251	1254	rBV	2191	3562	0.87%	0.069%
38	10.487	1262	1268	1276	rVB	60447	110887	27.21%	2.162%
39	10.628	1286	1292	1303	rBV3	41062	83646	20.53%	1.631%
40	10.869	1322	1333	1346	rVB3	118774	294528	72.27%	5.743%
41	11.010	1350	1357	1365	rVV3	19391	38261	9.39%	0.746%
42	11.086	1369	1370	1373	rBV2	1914	1727	0.42%	0.034%
43	11.574	1449	1453	1456	rVV2	2163	3043	0.75%	0.059%
44	11.674	1467	1470	1472	rBV	1620	1970	0.48%	0.038%
45	11.773	1484	1487	1493	rVB3	2815	3792	0.93%	0.074%
46	11.885	1500	1506	1510	rBV3	7602	12254	3.01%	0.239%
47	12.191	1553	1558	1560	rBV2	1823	2810	0.69%	0.055%
48	12.279	1566	1573	1578	rBV3	15912	21958	5.39%	0.428%
49	12.408	1593	1595	1599	rBV	1250	1683	0.41%	0.033%
50	12.525	1611	1615	1621	rVB2	10701	16277	3.99%	0.317%
51	12.749	1646	1653	1657	rBV2	6871	11092	2.72%	0.216%
52	12.819	1660	1665	1667	rVB3	2062	2358	0.58%	0.046%
53	13.083	1705	1710	1714	rVV4	3131	4901	1.20%	0.096%
54	13.172	1723	1725	1728	rBV2	1301	1588	0.39%	0.031%
55	13.289	1742	1745	1746	rBV2	1941	1538	0.38%	0.030%
56	13.365	1756	1758	1761	rBV2	1347	1724	0.42%	0.034%
57	13.512	1778	1783	1790	rBV4	19247	30935	7.59%	0.603%
58	13.583	1790	1795	1802	rVV3	24186	42731	10.49%	0.833%
59	13.853	1838	1841	1843	rBV2	5605	6837	1.68%	0.133%
60	14.017	1865	1869	1875	rVV3	10645	18742	4.60%	0.365%
61	14.094	1875	1882	1890	rVB	120986	195654	48.01%	3.815%
62	14.170	1890	1895	1899	rBV3	5987	9373	2.30%	0.183%
63	14.205	1899	1901	1904	rBV3	2322	2913	0.71%	0.057%
64	14.252	1906	1909	1913	rBV3	6467	7551	1.85%	0.147%
65	14.393	1927	1933	1941	rVB	132862	208795	51.23%	4.071%
66	14.529	1952	1956	1959	rBV2	2753	3359	0.82%	0.065%
67	14.581	1961	1965	1970	rVB2	19013	25925	6.36%	0.505%
68	14.699	1973	1985	1991	rBV	207852	336575	82.59%	6.562%
69	14.775	1995	1998	2000	rBV3	2660	3114	0.76%	0.061%
70	14.852	2007	2011	2012	rBV3	3589	3376	0.83%	0.066%
71	14.893	2013	2018	2029	rVV2	52068	98773	24.24%	1.926%
72	15.040	2040	2043	2047	rVB5	4344	5918	1.45%	0.115%
73	15.098	2047	2053	2059	rBV5	12631	24125	5.92%	0.470%
74	15.316	2086	2090	2093	rBV3	5787	8413	2.06%	0.164%
75	15.363	2094	2098	2104	rVB3	8857	12795	3.14%	0.249%
76	15.480	2114	2118	2122	rBV6	7224	11693	2.87%	0.228%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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

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 Title : SVOA CALIBRATION

77	15.533	2123	2127	2134	rVB2	25019	32683	8.02%	0.637%
78	15.586	2134	2136	2137	rBV2	1928	1799	0.44%	0.035%
79	15.692	2149	2154	2160	rVB2	174269	268867	65.97%	5.242%
80	15.809	2169	2174	2181	rVB3	51388	79080	19.40%	1.542%
81	16.021	2207	2210	2211	rBV	3186	3234	0.79%	0.063%
82	16.132	2226	2229	2230	rBV2	6586	6778	1.66%	0.132%
83	16.397	2270	2274	2283	rBV2	26267	44654	10.96%	0.871%
84	16.667	2318	2320	2322	rBV	1772	1909	0.47%	0.037%
85	16.720	2327	2329	2330	rBV	2470	1509	0.37%	0.029%
86	16.955	2364	2369	2370	rBV2	2902	3813	0.94%	0.074%
87	16.978	2370	2373	2376	rBV4	6489	8480	2.08%	0.165%
88	17.102	2391	2394	2399	rBV5	9308	15492	3.80%	0.302%
89	17.190	2405	2409	2413	rVB4	29053	42333	10.39%	0.825%
90	17.448	2447	2453	2458	rBV	266681	404297	99.21%	7.883%
91	17.548	2465	2470	2477	rVB2	193895	287938	70.65%	5.614%
92	17.930	2532	2535	2541	rVB2	28930	37301	9.15%	0.727%
93	18.106	2562	2565	2568	rBV3	14158	15452	3.79%	0.301%
94	18.329	2599	2603	2607	rBV2	71676	100605	24.69%	1.962%
95	18.511	2631	2634	2638	rVB3	19961	24638	6.05%	0.480%
96	18.623	2649	2653	2658	rBV2	31406	43446	10.66%	0.847%
97	18.823	2684	2687	2692	rVB5	20935	26487	6.50%	0.516%
98	19.246	2755	2759	2762	rBV4	32825	59285	14.55%	1.156%
99	19.833	2854	2859	2870	rVB	227455	381872	93.70%	7.445%
100	21.731	3177	3182	3188	rVB	241638	407532	100.00%	7.946%

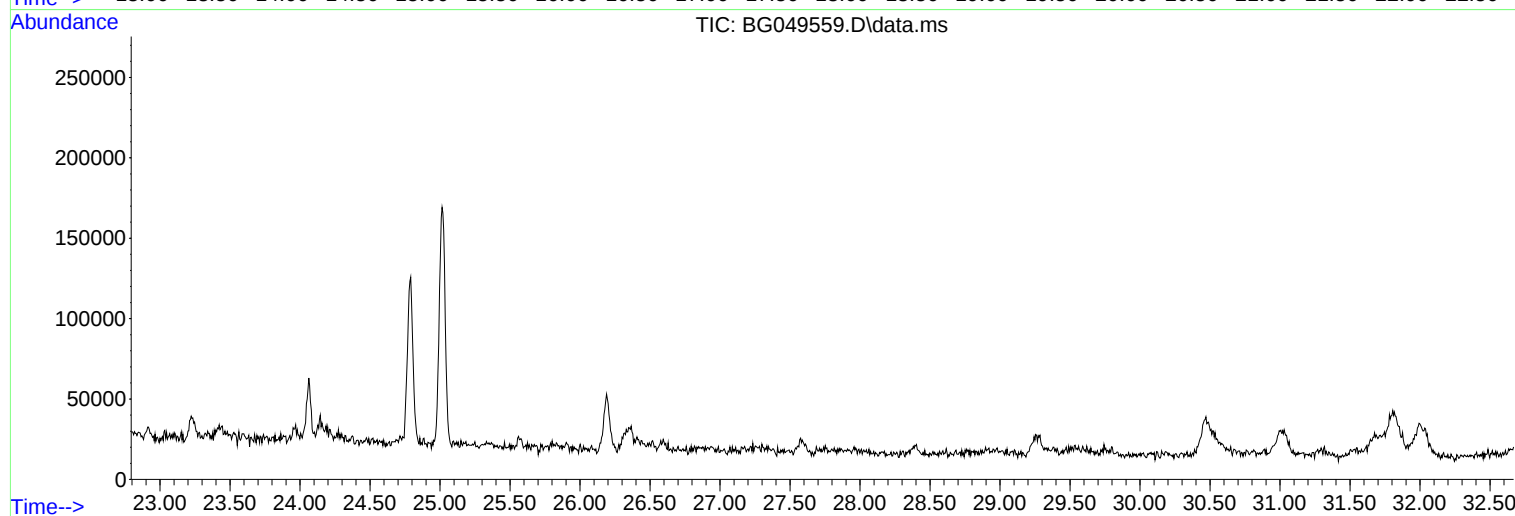
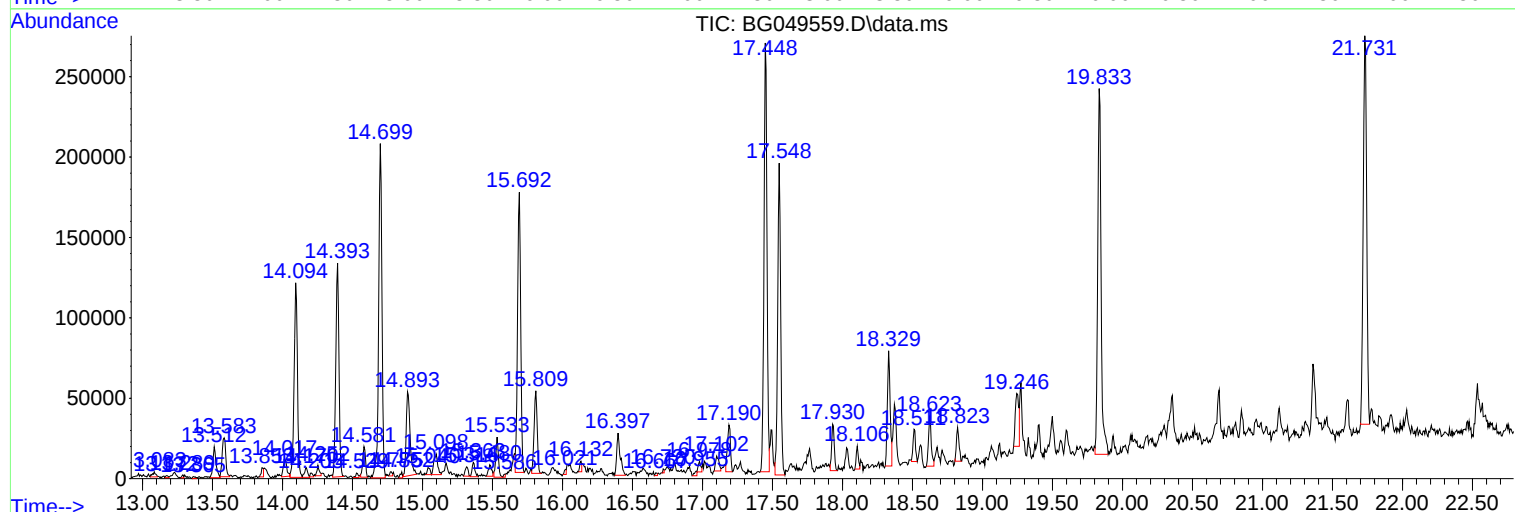
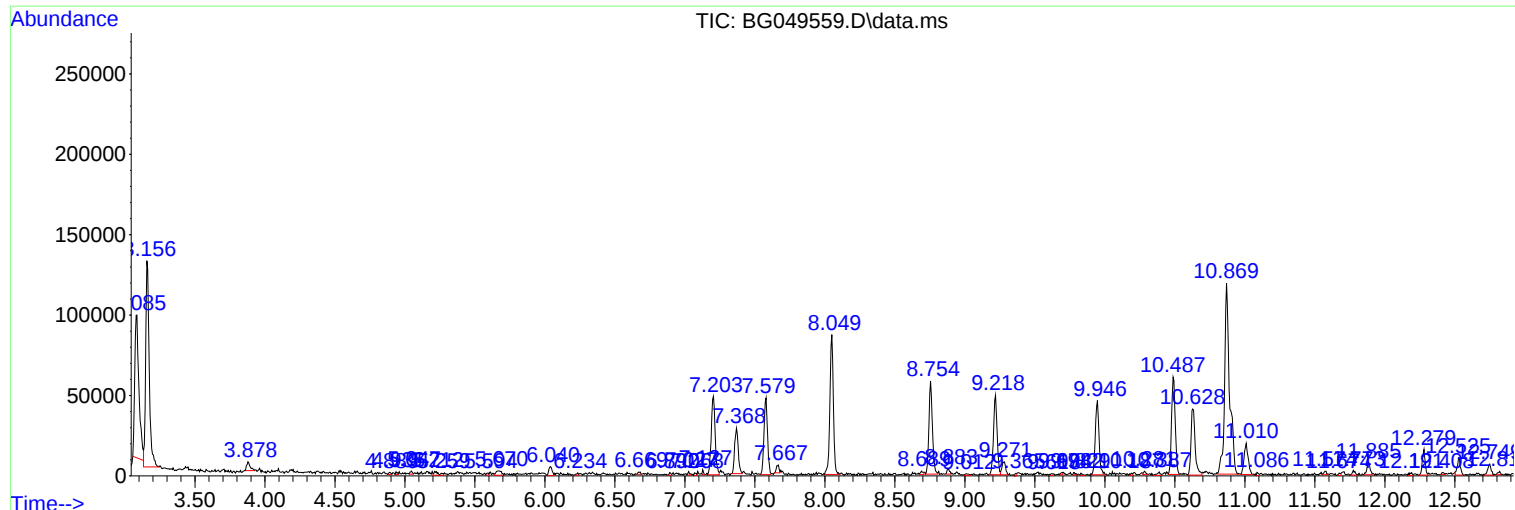
Sum of corrected areas: 5128908

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Instrument :
BNA_G
ClientSampled :
C0068

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
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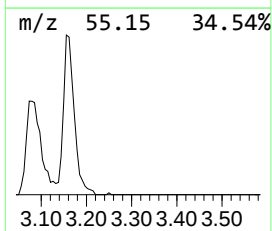
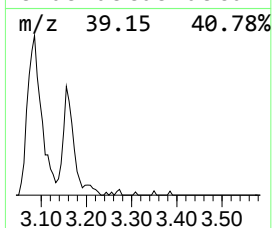
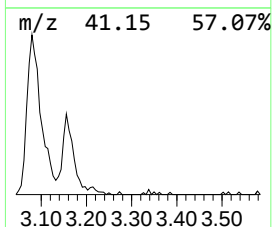
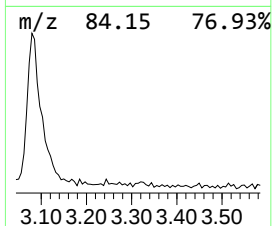
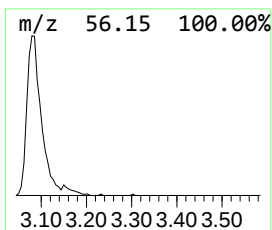
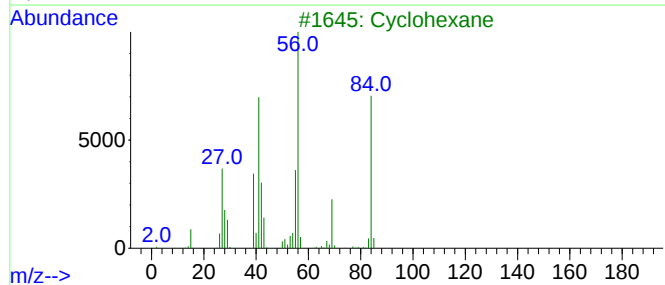
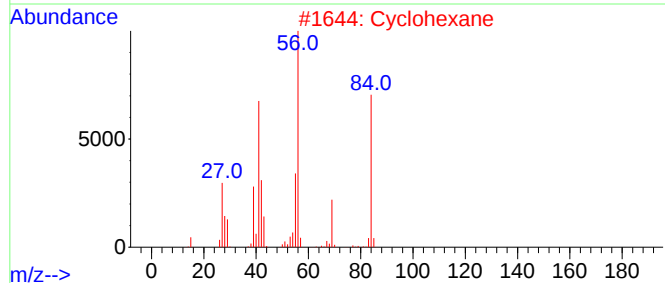
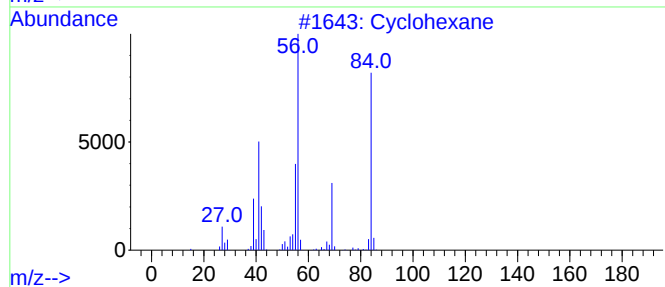
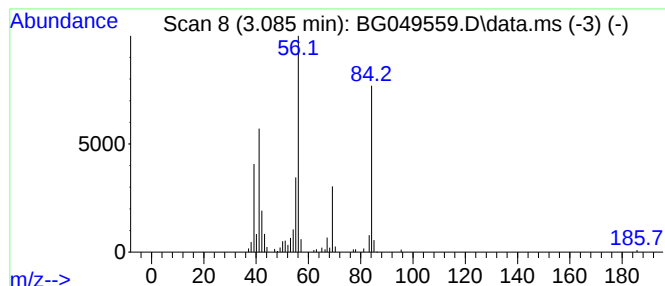
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.085 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.085	24.68 ng/ul	186903	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	94
2			Cyclohexane	84	C6H12	000110-82-7	81
3			Cyclohexane	84	C6H12	000110-82-7	81
4			2-Amino-oxazole	84	C3H4N2O	004570-45-0	72
5			Cyclohexane	84	C6H12	000110-82-7	68



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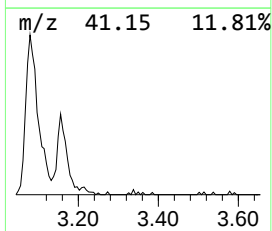
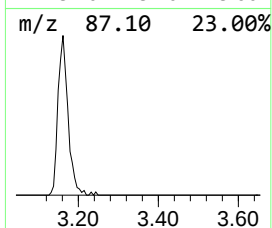
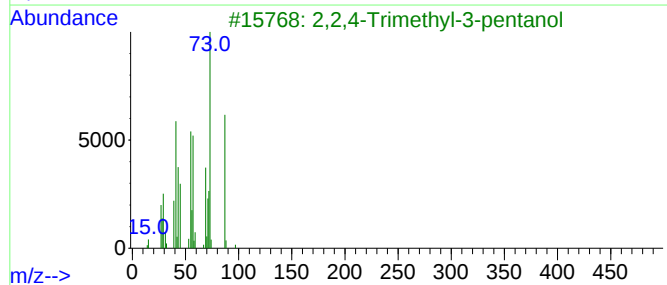
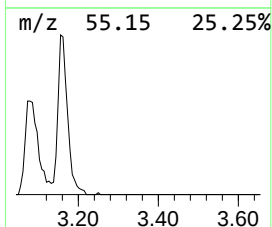
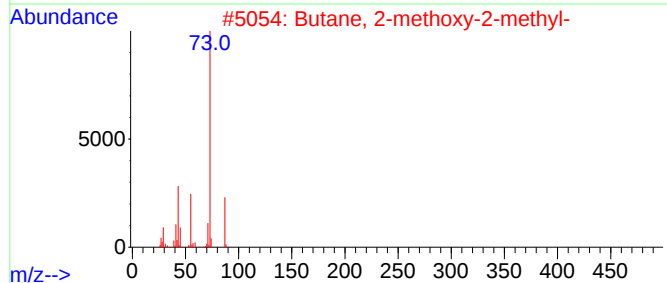
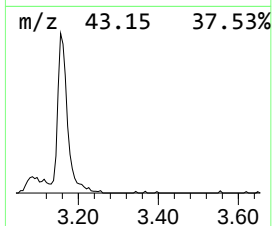
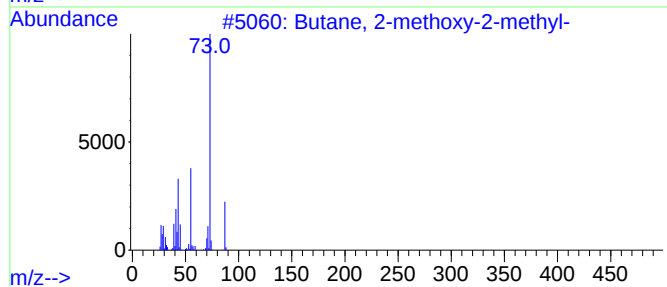
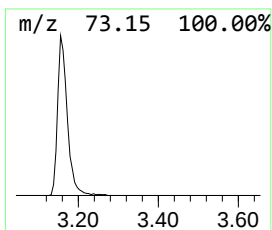
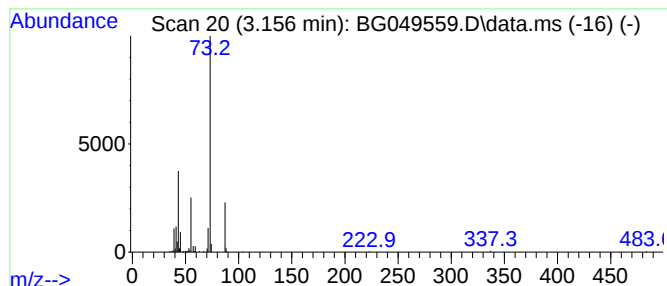
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.156	29.39 ng/ul	222609	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	33



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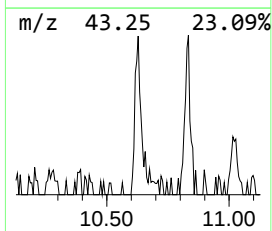
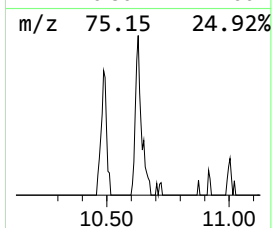
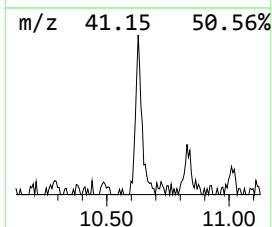
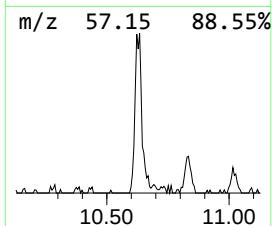
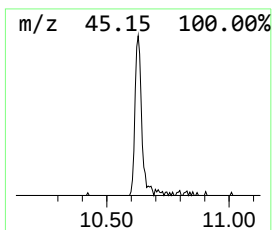
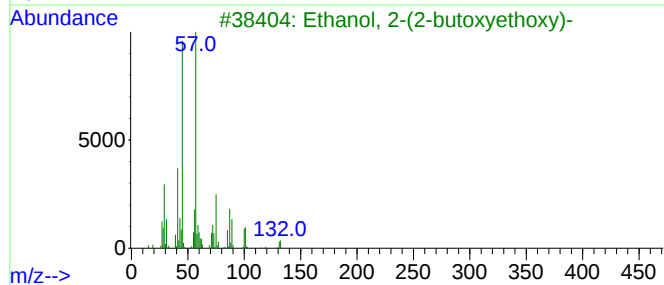
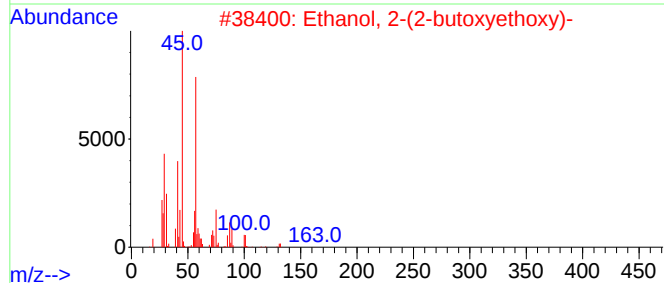
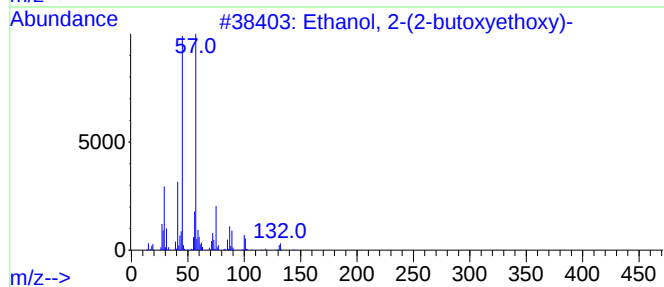
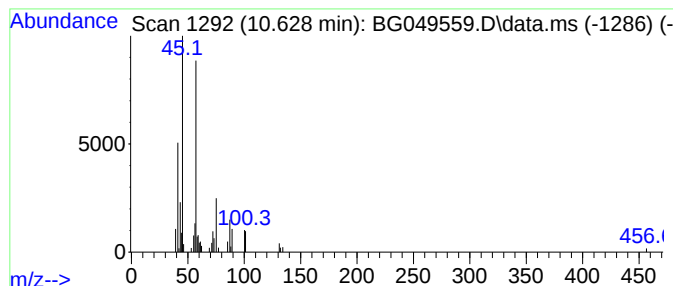
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.628	5.68 ng/ul	83646	Naphthalene-d8	10.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
Operator : CG/JU
Sample : M3123-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

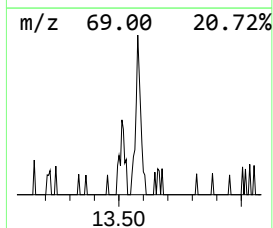
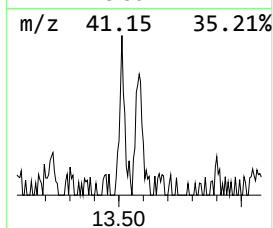
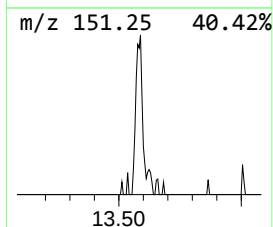
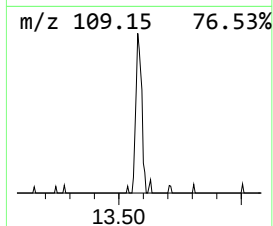
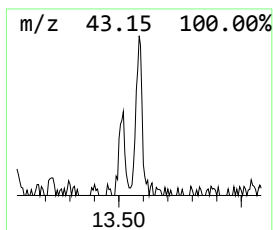
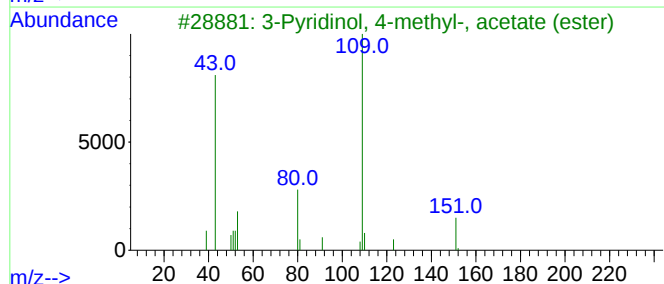
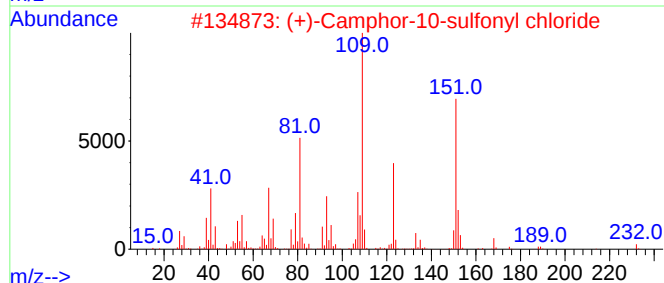
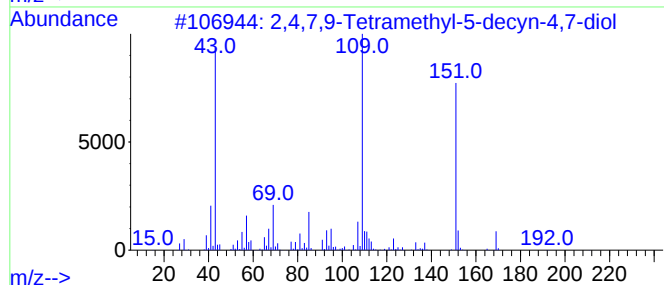
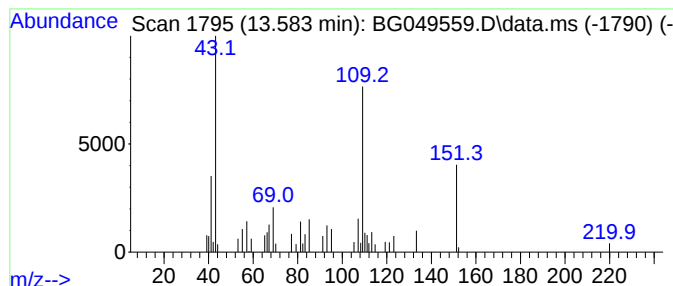
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.583	2.54 ng/ul	42731	Acenaphthene-d10	14.699	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	72
2	(+)-Camphor-10-sulfonyl chloride	250	C10H15ClO3S	021286-54-4	59
3	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	50
4	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	47
5	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
Operator : CG/JU
Sample : M3123-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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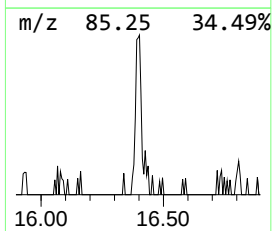
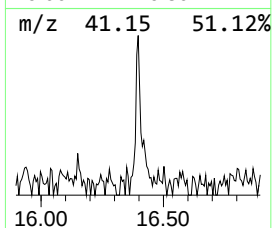
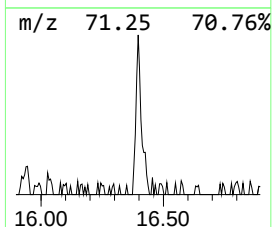
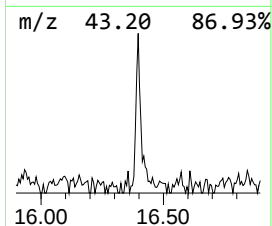
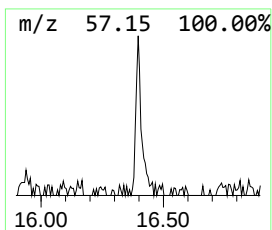
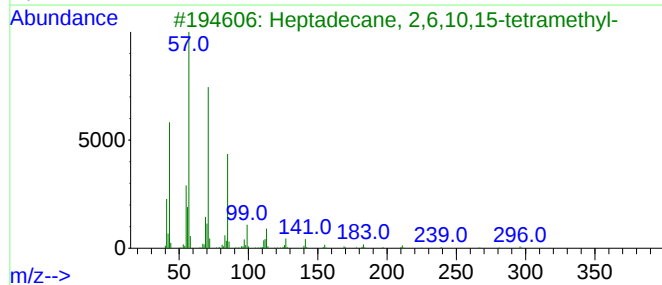
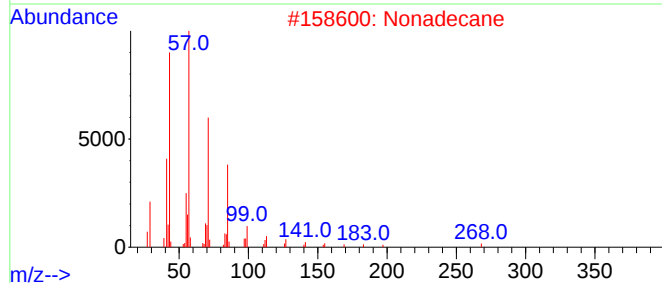
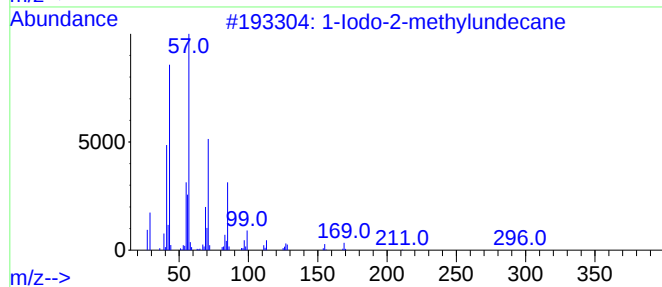
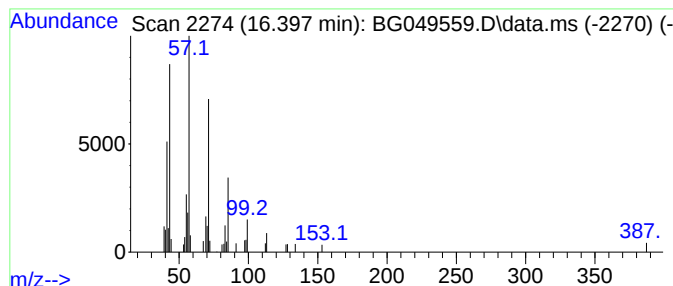
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Iodo-2-methylundecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.397	2.21 ng/ul	44654	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	80
2			Nonadecane	268	C19H40	000629-92-5	74
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
5			Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
Operator : CG/JU
Sample : M3123-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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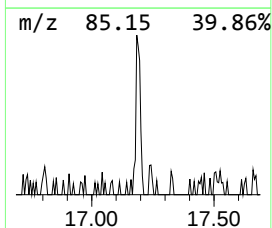
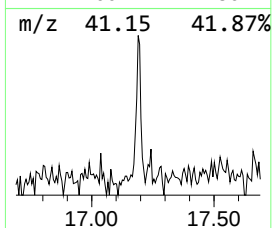
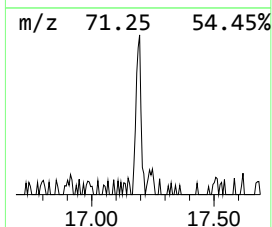
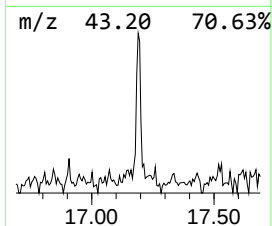
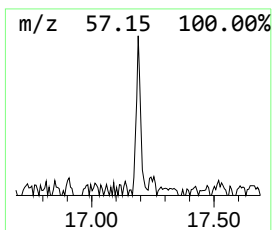
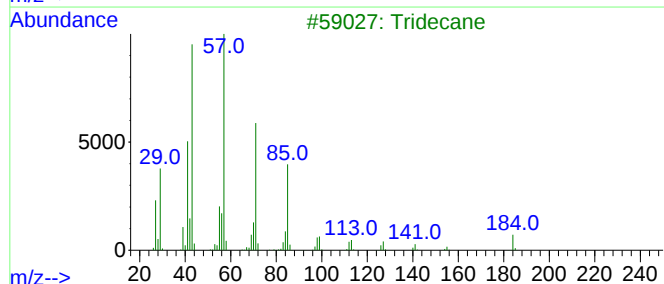
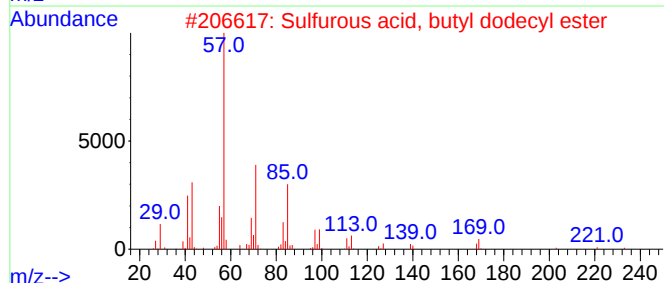
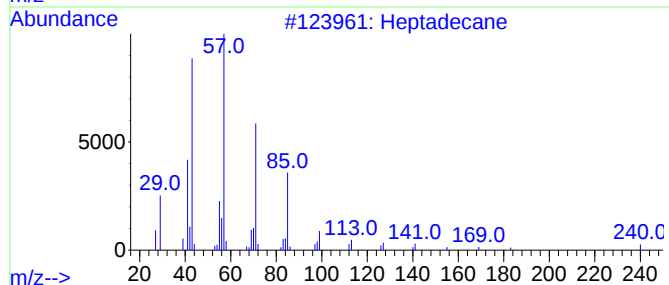
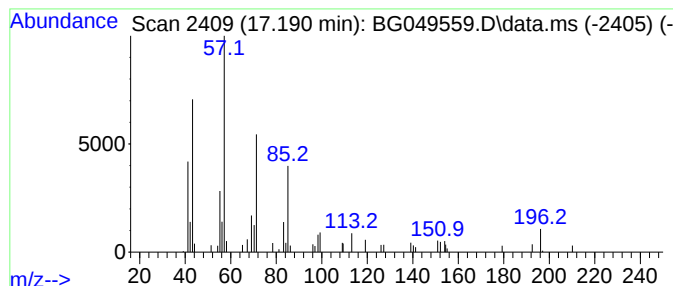
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.190	2.09 ng/ul	42333	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	64
2			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	64
3			Tridecane	184	C13H28	000629-50-5	59
4			Tetratetracontane	619	C44H90	007098-22-8	59
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
Operator : CG/JU
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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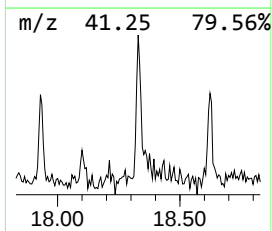
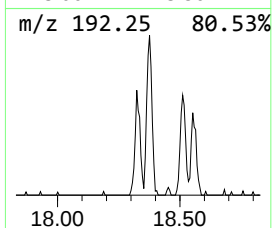
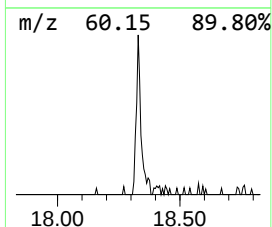
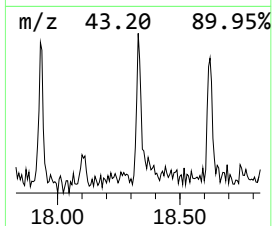
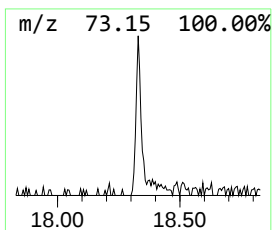
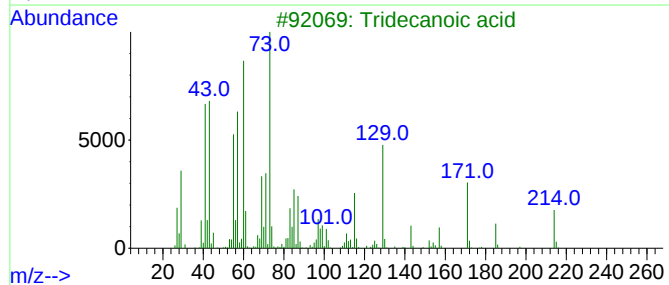
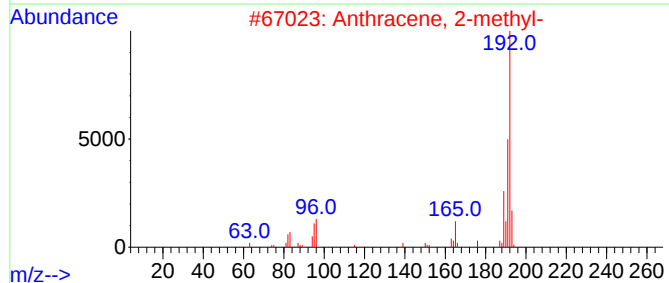
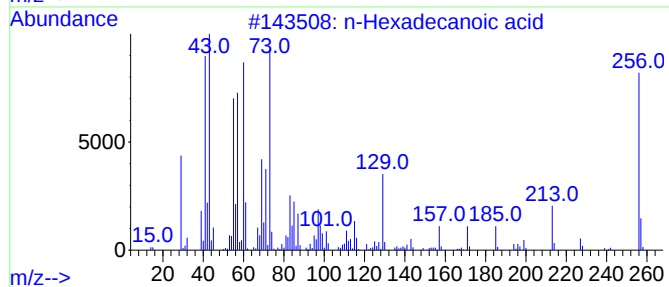
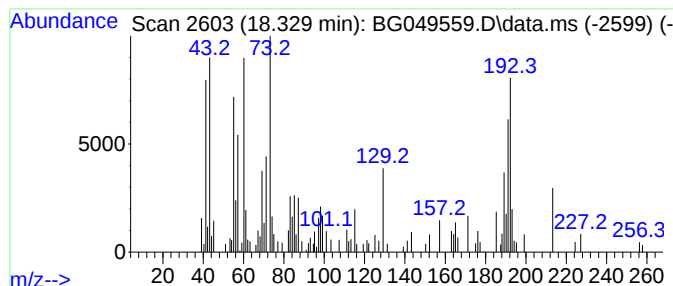
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	4.98 ng/ul	100605	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	52
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	43
3			Tridecanoic acid	214	C13H26O2	000638-53-9	38
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	38
5			8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
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Sample : M3123-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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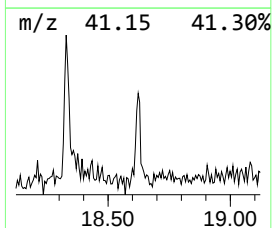
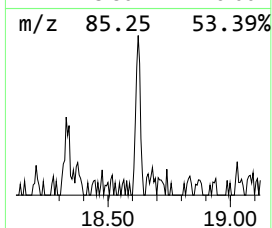
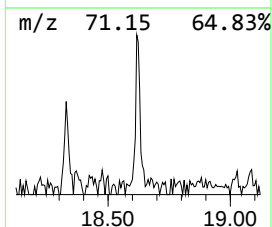
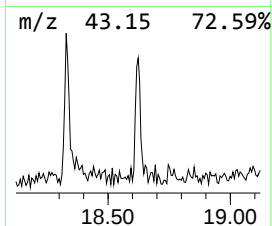
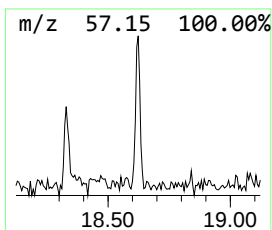
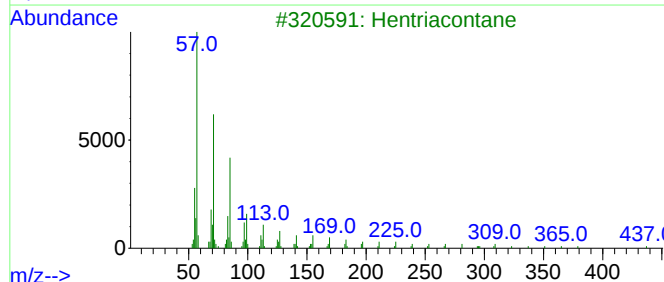
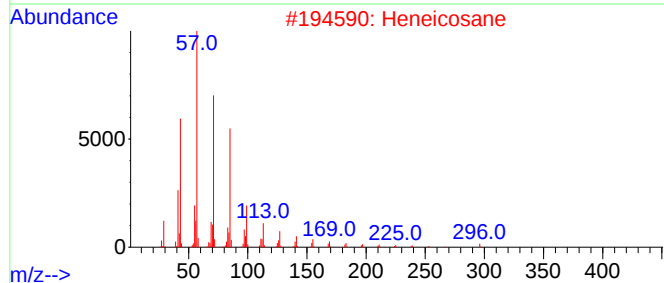
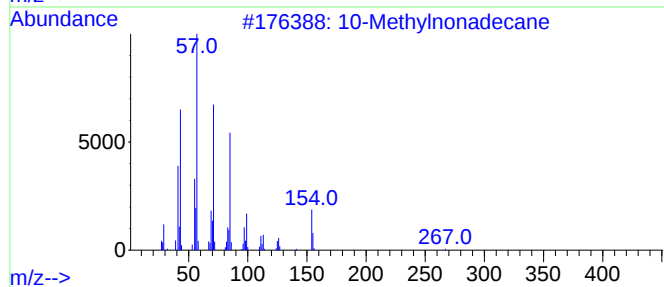
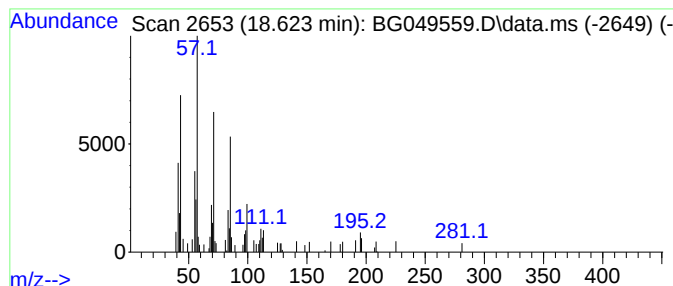
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.623	2.15 ng/ul	43446	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane			282	C20H42	056862-62-5	72
2	Heneicosane			296	C21H44	000629-94-7	64
3	Hentriacontane			437	C31H64	000630-04-6	64
4	Hexadecane			226	C16H34	000544-76-3	58
5	Heptadecane			240	C17H36	000629-78-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
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Misc :
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Instrument :
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ClientSampleId :
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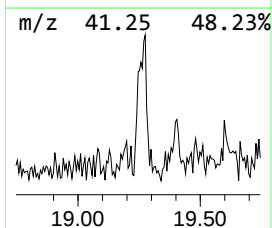
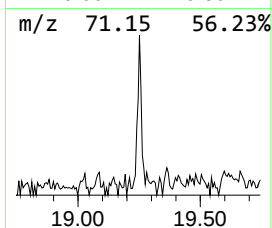
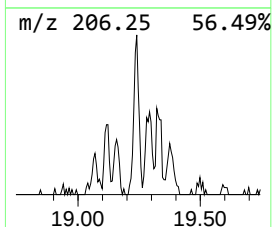
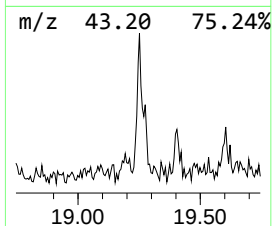
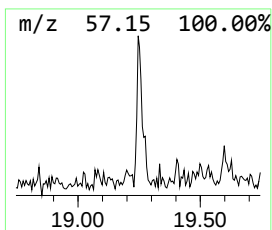
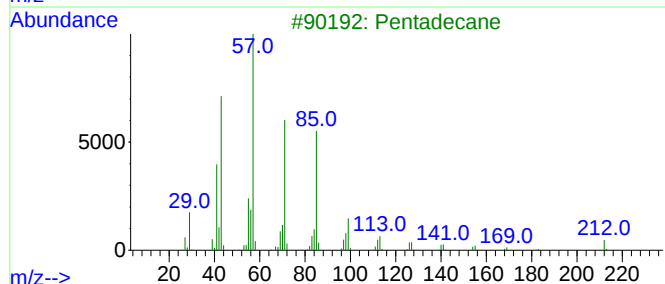
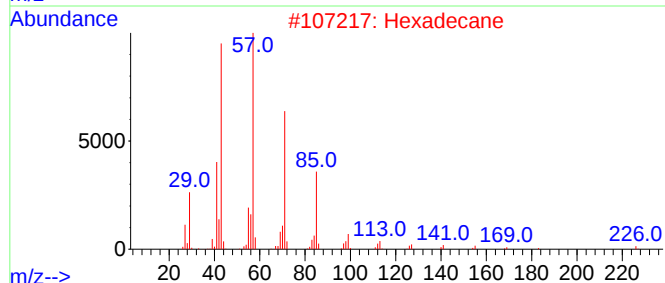
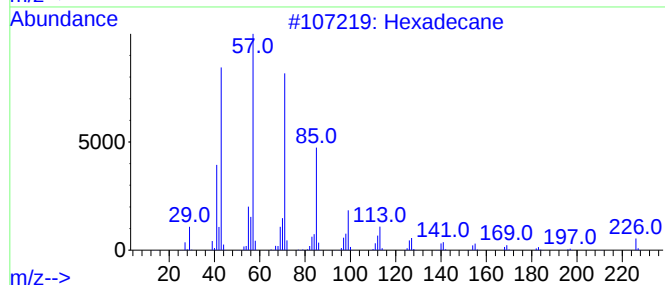
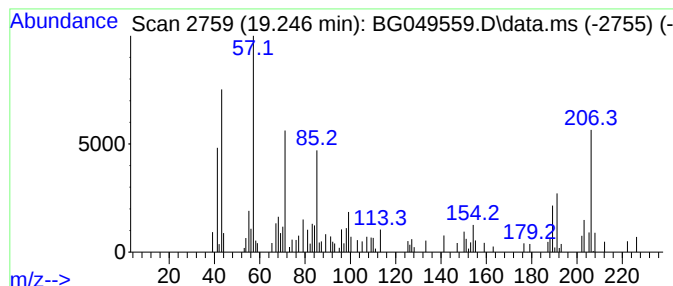
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.246	2.93 ng/ul	59285	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			Hexadecane	226	C16H34	000544-76-3	49
3			Pentadecane	212	C15H32	000629-62-9	42
4			Hexadecane	226	C16H34	000544-76-3	42
5			Pentadecane	212	C15H32	000629-62-9	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049559.D
Acq On : 29 Jul 2021 14:32
Operator : CG/JU
Sample : M3123-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Butane, 2-metho...	3.156	29.4	ng/ul	222609	1	8.049	151467	20.0
Ethanol, 2-(2-b...	10.628	5.7	ng/ul	83646	2	10.869	294528	20.0
2,4,7,9-Tetrame...	13.583	2.5	ng/ul	42731	3	14.699	336575	20.0
1-Iodo-2-methyl...	16.397	2.2	ng/ul	44654	4	17.448	404297	20.0
(DEL) Alkane: S...	17.190	2.1	ng/ul	42333	4	17.448	404297	20.0
n-Hexadecanoic ...	18.329	5.0	ng/ul	100605	4	17.448	404297	20.0
(DEL) Alkane: S...	18.623	2.1	ng/ul	43446	4	17.448	404297	20.0
(DEL) Alkane: S...	19.246	2.9	ng/ul	59285	4	17.448	404297	20.0