

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036561.D
 Acq On : 16 Sep 2022 14:10
 Operator : CG/JU
 Sample : N4601-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5753

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_M\Methods\SFAM-EPA-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036561.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	30	rVB	12119659	16421533	100.00%	12.639%
2	6.022	493	499	508	rVV3	114307	224753	1.37%	0.173%
3	7.169	688	694	703	rVV	997544	1670623	10.17%	1.286%
4	7.351	719	725	731	rVV	857156	1351894	8.23%	1.041%
5	7.551	749	759	767	rBV	943554	1525124	9.29%	1.174%
6	8.022	832	839	846	rBV	1208846	2005147	12.21%	1.543%
7	8.722	945	958	967	rBV	1038747	1722213	10.49%	1.326%
8	8.851	974	980	987	rVV3	66795	166077	1.01%	0.128%
9	9.186	1031	1037	1043	rVV	945282	1498752	9.13%	1.154%
10	9.263	1046	1050	1059	rVB	105042	194948	1.19%	0.150%
11	9.910	1151	1160	1166	rBV	642032	1048266	6.38%	0.807%
12	10.445	1243	1251	1260	rVB	1148470	1936389	11.79%	1.490%
13	10.610	1272	1279	1286	rBV	807129	1238596	7.54%	0.953%
14	10.839	1308	1318	1323	rBV	1783521	3212621	19.56%	2.473%
15	10.886	1323	1326	1333	rVB	196345	299316	1.82%	0.230%
16	10.969	1333	1340	1345	rBV	650170	1124513	6.85%	0.866%
17	11.869	1487	1493	1501	rVV	210718	343280	2.09%	0.264%
18	12.257	1554	1559	1570	rBV	175735	299035	1.82%	0.230%
19	12.486	1592	1598	1605	rVB	365276	612257	3.73%	0.471%
20	12.704	1629	1635	1644	rVV	335961	527123	3.21%	0.406%
21	13.204	1716	1720	1725	rVV	153111	211301	1.29%	0.163%
22	13.486	1762	1768	1774	rBV	319697	463808	2.82%	0.357%
23	13.551	1774	1779	1784	rBV	123512	197709	1.20%	0.152%
24	13.815	1819	1824	1826	rBV2	123895	207835	1.27%	0.160%
25	13.974	1846	1851	1856	rVV	338428	586089	3.57%	0.451%
26	14.027	1856	1860	1861	rVV	288654	372976	2.27%	0.287%
27	14.057	1861	1865	1875	rVV	2318220	3238556	19.72%	2.493%
28	14.139	1875	1879	1883	rVV	264895	349090	2.13%	0.269%
29	14.210	1883	1891	1894	rVV2	197306	353430	2.15%	0.272%
30	14.345	1907	1914	1930	rVB	2625401	4111069	25.03%	3.164%
31	14.551	1944	1949	1956	rVB	264868	343444	2.09%	0.264%
32	14.645	1956	1965	1972	rBV	2862790	4481473	27.29%	3.449%
33	14.821	1989	1995	2000	rBV	506284	747835	4.55%	0.576%
34	15.045	2028	2033	2039	rVV	301276	487731	2.97%	0.375%
35	15.121	2039	2046	2050	rVV	142212	228409	1.39%	0.176%
36	15.262	2063	2070	2075	rVV2	142293	260860	1.59%	0.201%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Title : SVOA CALIBRATION

37	15.315	2075	2079	2083	rVB	143408	187885	1.14%	0.145%
38	15.433	2092	2099	2107	rBV2	219737	490194	2.99%	0.377%
39	15.498	2107	2110	2116	rVB	266315	319230	1.94%	0.246%
40	15.633	2126	2133	2138	rVV	3378520	4987915	30.37%	3.839%
41	15.680	2138	2141	2145	rVB	220702	284260	1.73%	0.219%
42	15.739	2145	2151	2156	rBV3	680082	950260	5.79%	0.731%
43	15.904	2175	2179	2184	rVB	201964	285520	1.74%	0.220%
44	15.986	2185	2193	2201	rVB	204017	394339	2.40%	0.304%
45	16.127	2212	2217	2225	rVB2	185693	433348	2.64%	0.334%
46	16.215	2225	2232	2240	rVB3	112797	254805	1.55%	0.196%
47	16.362	2252	2257	2258	rBV2	402131	547599	3.33%	0.421%
48	16.380	2258	2260	2270	rVB	645342	872082	5.31%	0.671%
49	16.921	2348	2352	2356	rBV3	168739	287388	1.75%	0.221%
50	17.039	2368	2372	2379	rBV4	123879	232779	1.42%	0.179%
51	17.115	2379	2385	2387	rVV2	129552	242006	1.47%	0.186%
52	17.151	2387	2391	2396	rVV	385389	582121	3.54%	0.448%
53	17.204	2396	2400	2410	rVB3	203054	356450	2.17%	0.274%
54	17.386	2424	2431	2435	rBV	3443082	5009033	30.50%	3.855%
55	17.427	2435	2438	2442	rVV	736850	991374	6.04%	0.763%
56	17.480	2442	2447	2457	rVB	3771246	5491752	33.44%	4.227%
57	17.568	2458	2462	2468	rVV3	99124	183238	1.12%	0.141%
58	17.692	2480	2483	2490	rVB2	167462	256460	1.56%	0.197%
59	17.892	2513	2517	2521	rVB	359366	442497	2.69%	0.341%
60	17.962	2526	2529	2533	rVB	131643	159198	0.97%	0.123%
61	18.062	2542	2546	2550	rVB	311295	370993	2.26%	0.286%
62	18.280	2575	2583	2585	rVV2	425364	1019016	6.21%	0.784%
63	18.309	2585	2588	2593	rVB	638291	871290	5.31%	0.671%
64	18.445	2606	2611	2615	rBV	452183	693842	4.23%	0.534%
65	18.486	2615	2618	2625	rVB	341261	485285	2.96%	0.374%
66	18.580	2630	2634	2642	rVB	397872	526198	3.20%	0.405%
67	18.756	2659	2664	2667	rBV	294390	398925	2.43%	0.307%
68	18.792	2667	2670	2677	rVB5	129086	207634	1.26%	0.160%
69	19.003	2702	2706	2710	rVV	129222	164070	1.00%	0.126%
70	19.050	2710	2714	2717	rVV	182921	251472	1.53%	0.194%
71	19.086	2718	2720	2724	rVB	167353	199678	1.22%	0.154%
72	19.168	2729	2734	2738	rBV	508426	773994	4.71%	0.596%
73	19.227	2738	2744	2747	rVV2	685411	1322769	8.06%	1.018%
74	19.262	2747	2750	2754	rVB	270731	329879	2.01%	0.254%
75	19.309	2754	2758	2764	rBV2	220206	458380	2.79%	0.353%
76	19.433	2774	2779	2786	rVB	1319393	1738745	10.59%	1.338%

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77	19.550	2793	2799	2802	rBV5	146957	258135	1.57%	0.199%
78	19.762	2830	2835	2848	rVV2	5048766	8452455	51.47%	6.506%
79	19.868	2849	2853	2859	rVB2	303389	477574	2.91%	0.368%
80	20.109	2890	2894	2906	rBV5	221552	651156	3.97%	0.501%
81	20.280	2921	2923	2926	rVB	197233	213285	1.30%	0.164%
82	20.315	2926	2929	2933	rVB2	396022	418542	2.55%	0.322%
83	20.439	2945	2950	2954	rBV2	239994	420273	2.56%	0.323%
84	20.768	3003	3006	3009	rVB	264448	254890	1.55%	0.196%
85	20.809	3010	3013	3020	rVB2	320041	362315	2.21%	0.279%
86	21.027	3046	3050	3057	rVB	538819	726079	4.42%	0.559%
87	21.197	3072	3079	3082	rBV3	248180	516987	3.15%	0.398%
88	21.268	3087	3091	3097	rVB3	886683	1281449	7.80%	0.986%
89	21.550	3132	3139	3143	rBV2	4563334	7340394	44.70%	5.650%
90	21.591	3143	3146	3156	rVB	1019378	1645507	10.02%	1.267%
91	21.721	3165	3168	3174	rBV4	419973	578429	3.52%	0.445%
92	22.197	3244	3249	3255	rBV4	812711	1392954	8.48%	1.072%
93	22.709	3332	3336	3343	rVB2	318272	503422	3.07%	0.387%
94	22.850	3356	3360	3368	rVB	1573278	2380508	14.50%	1.832%
95	23.250	3423	3428	3430	rBV	785647	1330272	8.10%	1.024%
96	23.780	3514	3518	3522	rBV	374985	620566	3.78%	0.478%
97	23.838	3522	3528	3534	rVV2	2834236	5196140	31.64%	3.999%
98	23.997	3548	3555	3562	rBV2	2498845	5049923	30.75%	3.887%
99	24.679	3667	3671	3678	rVB	263559	486031	2.96%	0.374%
100	27.532	4152	4156	4180	rVB7	391924	1247425	7.60%	0.960%

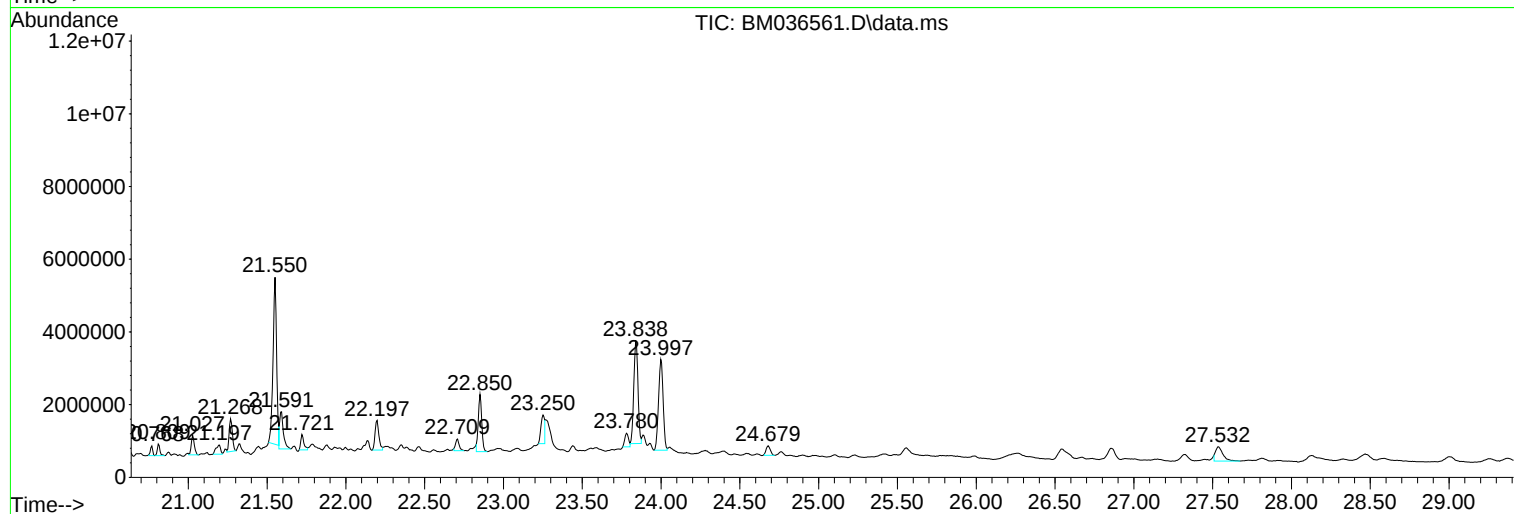
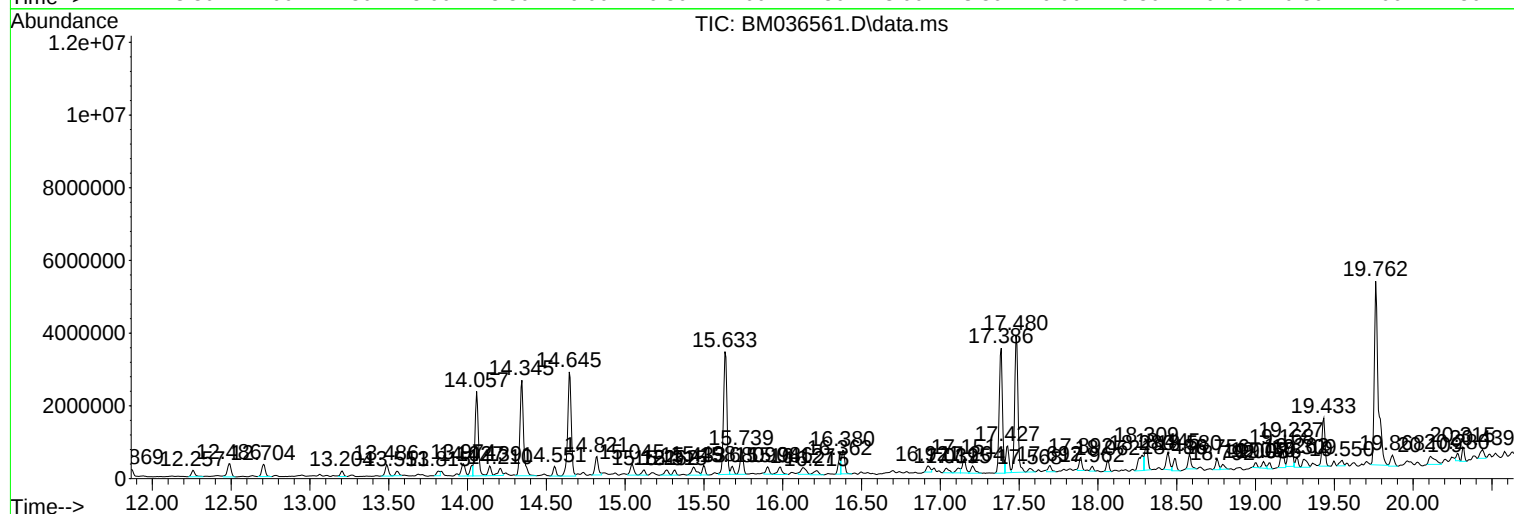
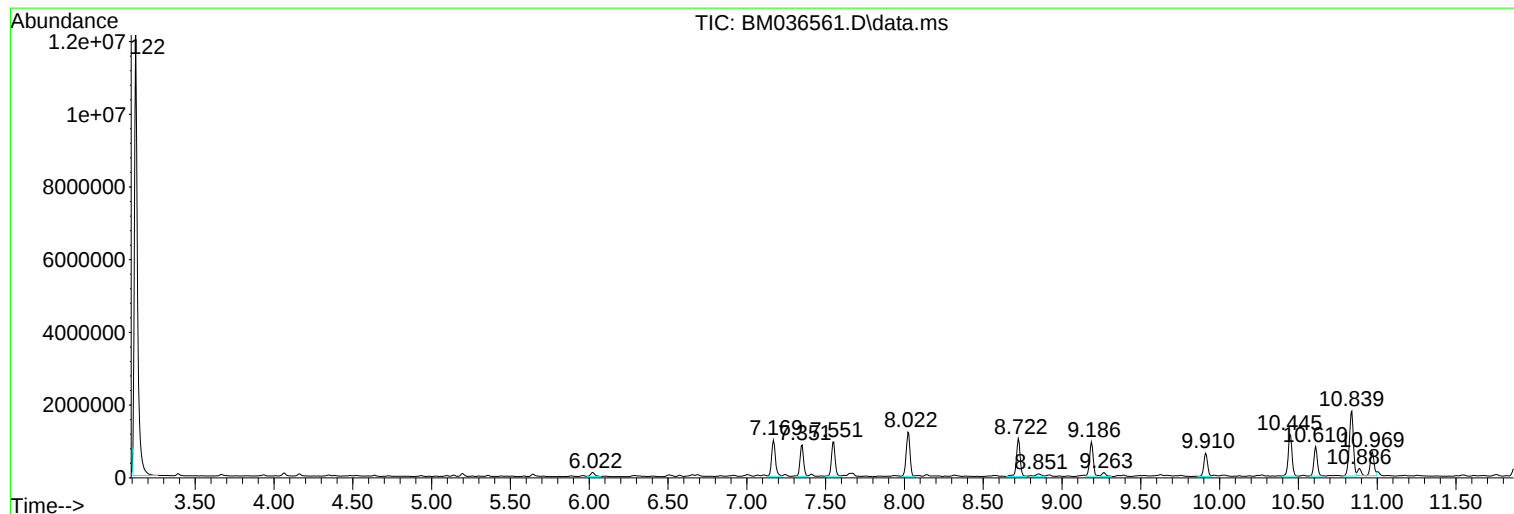
Sum of corrected areas: 129922689

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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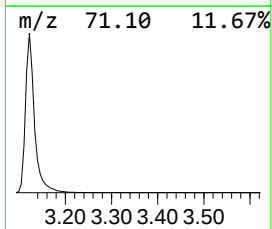
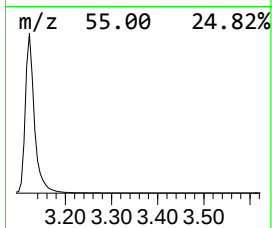
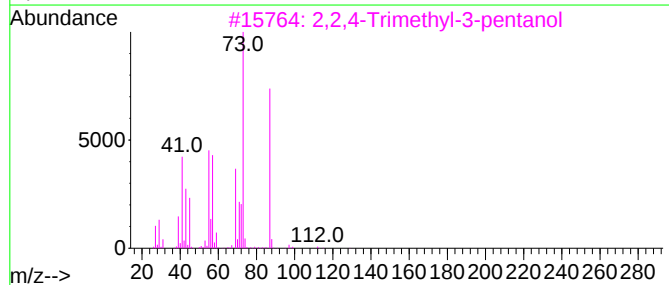
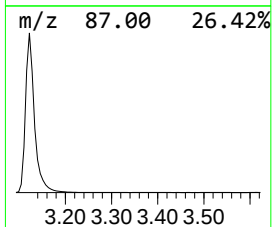
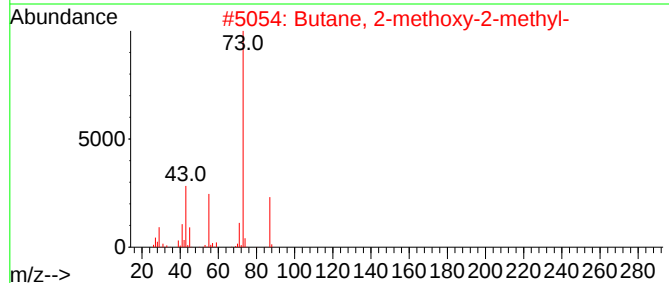
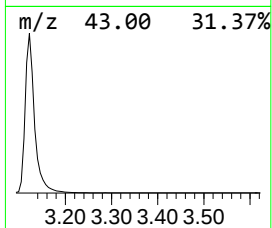
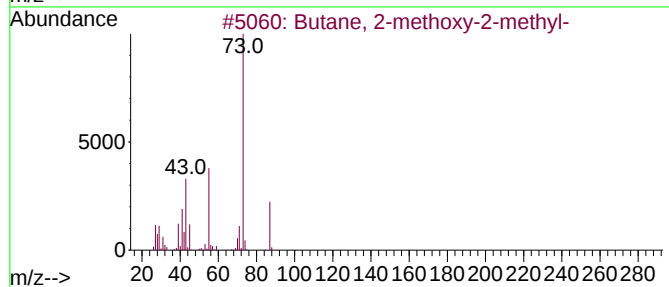
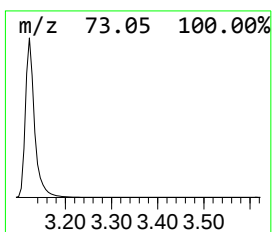
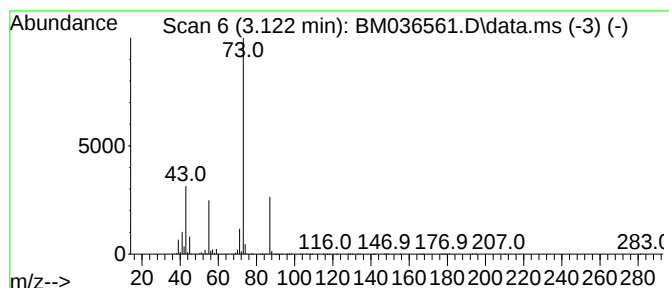
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	163.79 ng/ul	16421500	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9		



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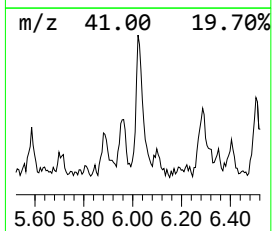
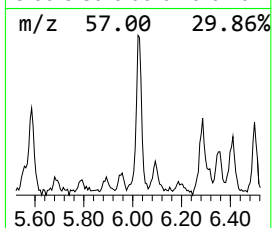
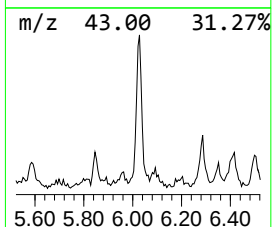
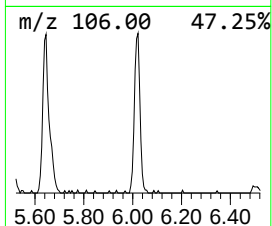
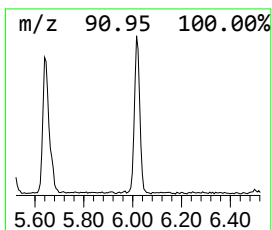
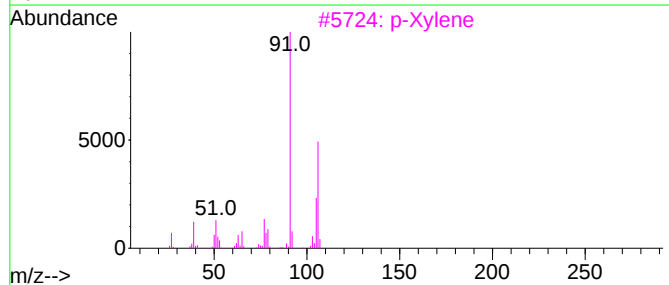
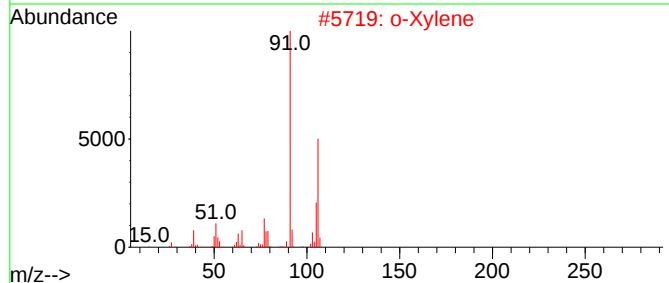
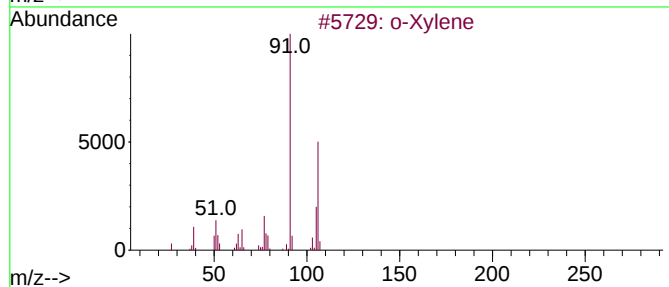
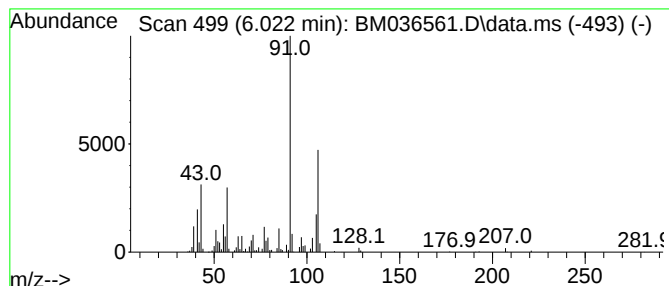
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 o-Xylene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.022	2.24 ng/ul	224753	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	94
2			o-Xylene	106	C8H10	000095-47-6	94
3			p-Xylene	106	C8H10	000106-42-3	94
4			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94



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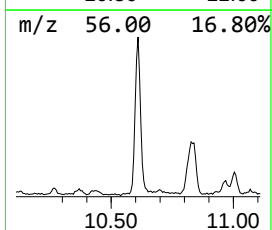
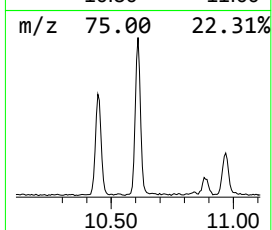
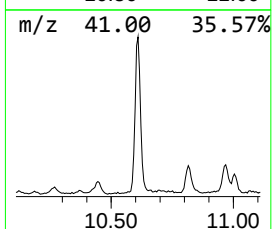
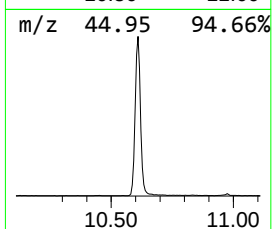
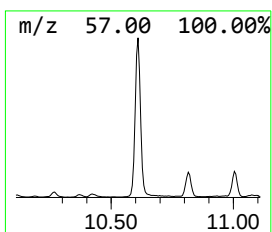
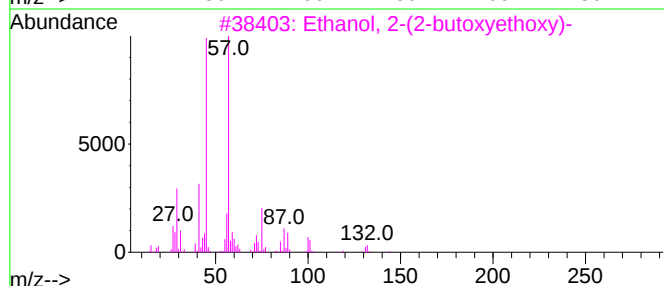
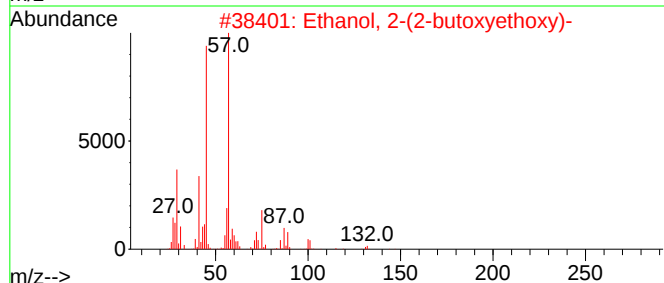
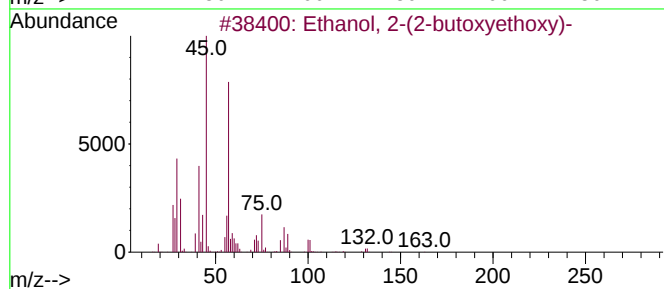
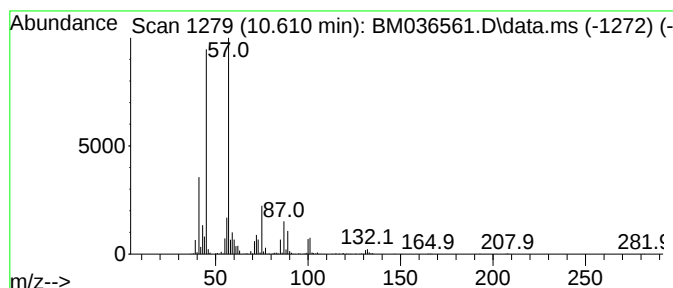
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.610	7.71 ng/ul	1238600	Naphthalene-d8	10.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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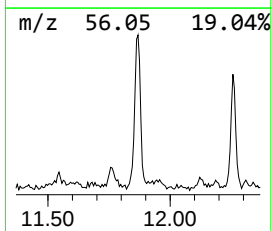
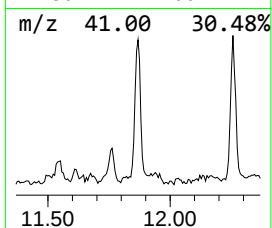
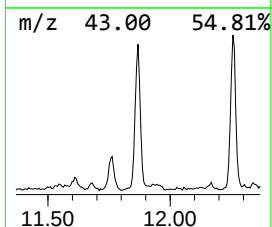
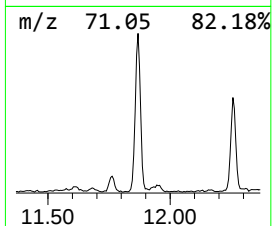
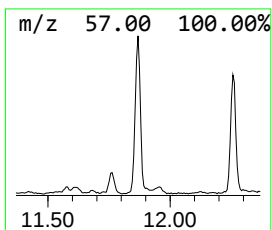
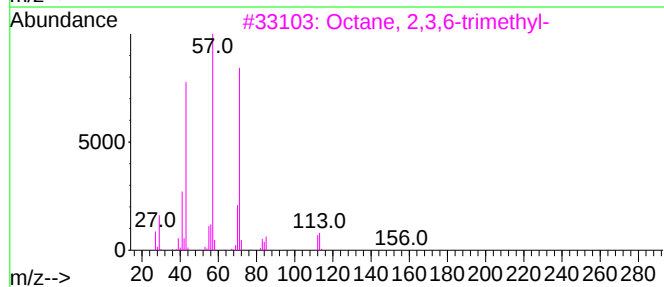
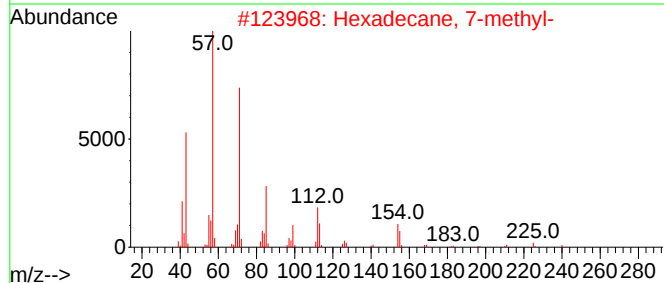
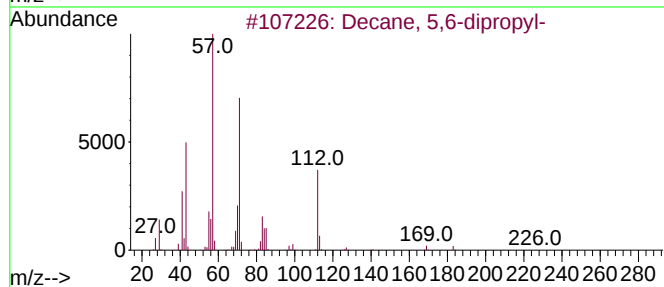
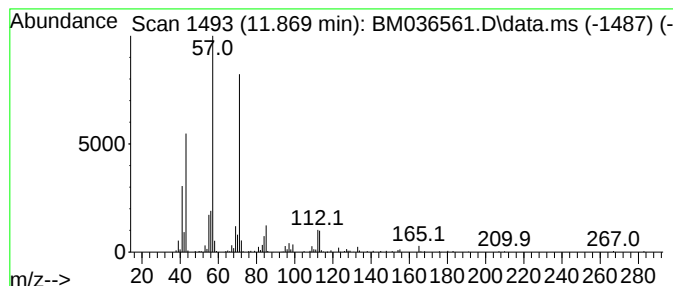
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.869	2.14 ng/ul	343280	Naphthalene-d8	10.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	78
2			Hexadecane, 7-methyl-	240	C17H36	026730-20-1	72
3			Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	72
4			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	64
5			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
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Sample : N4601-12
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Instrument :
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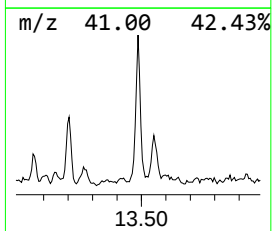
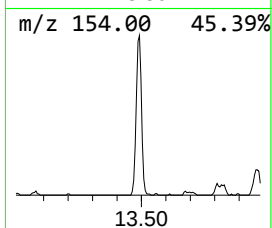
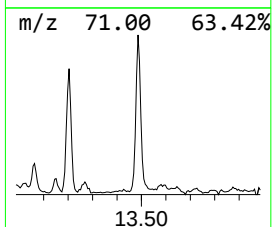
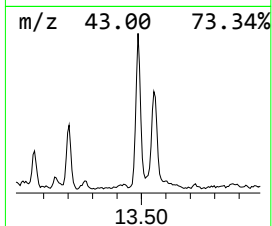
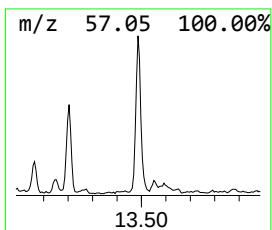
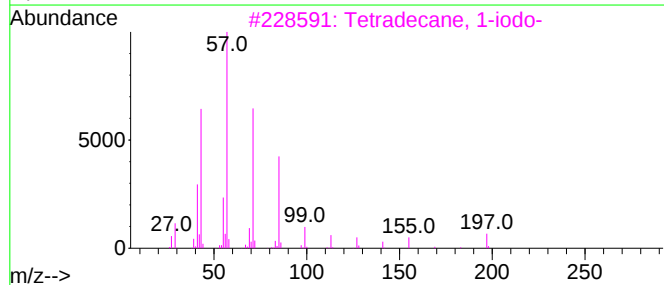
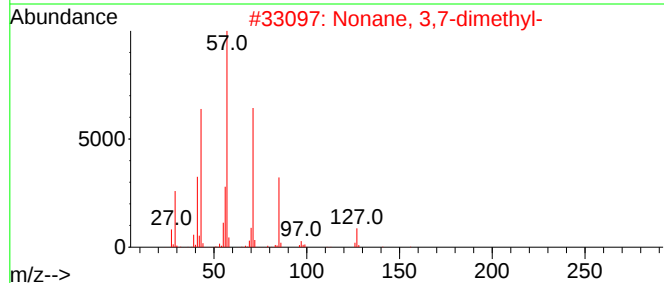
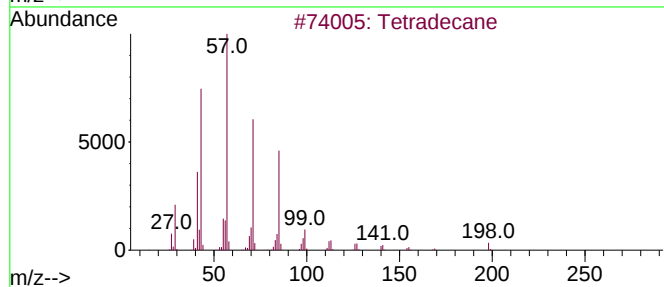
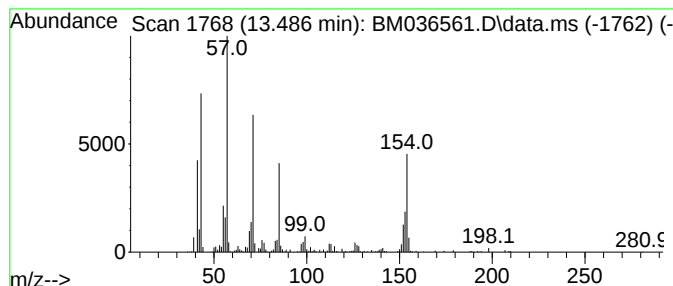
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.486	2.07 ng/ul	463808	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	91
2			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	49
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	46
4			Tridecane	184	C13H28	000629-50-5	46
5			Tridecane, 7-propyl-	226	C16H34	055045-09-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

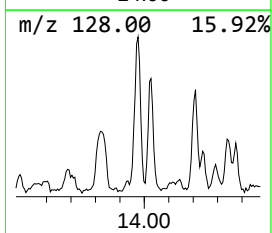
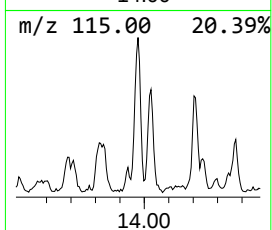
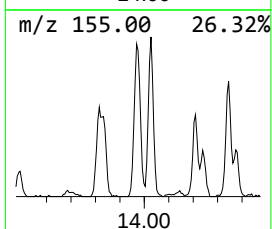
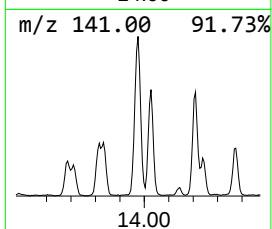
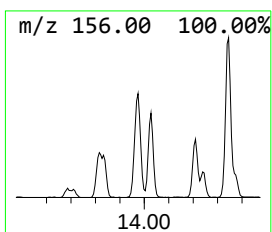
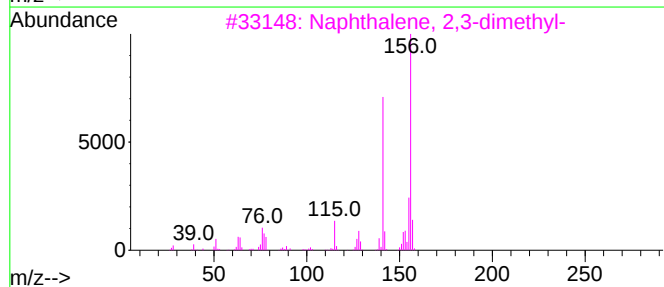
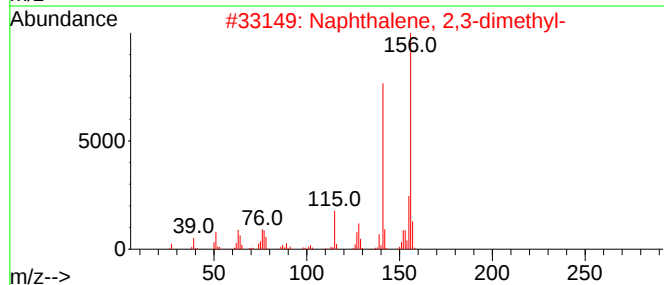
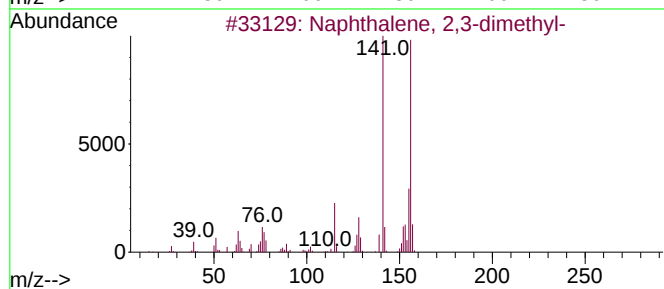
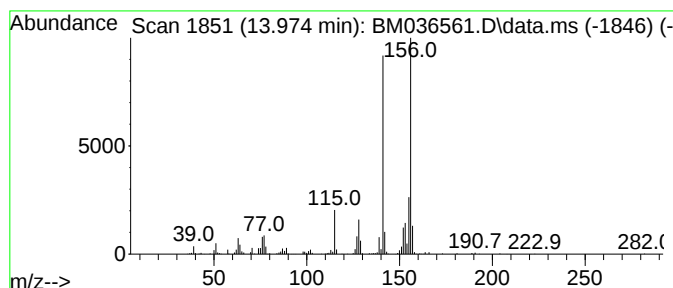
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.974	2.62 ng/ul	586089	Acenaphthene-d10		14.645	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	98
2	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97
3	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97
4	Naphthalene, 2,7-dimethyl-		156	C12H12	000582-16-1	97
5	Naphthalene, 1,6-dimethyl-		156	C12H12	000575-43-9	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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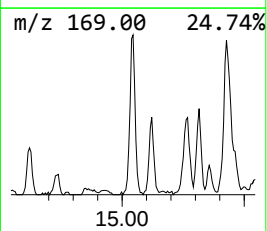
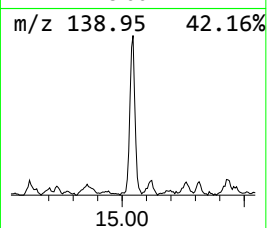
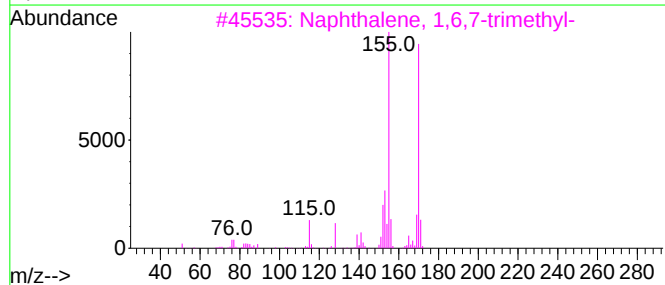
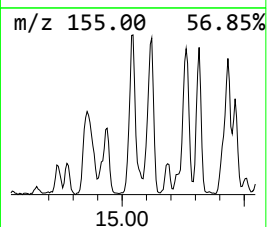
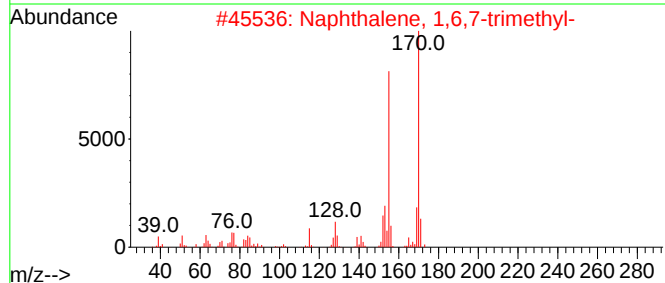
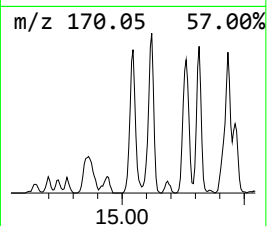
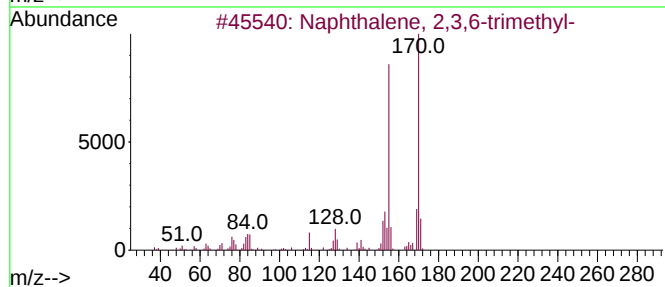
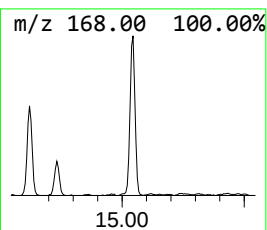
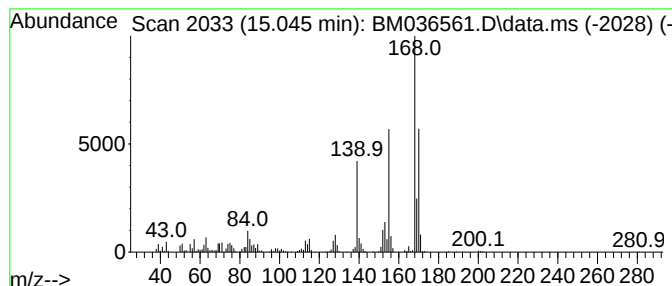
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene, 2,3,6-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.045	2.18 ng/ul	487731	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	83
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	55
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	52
5			Dibenzofuran	168	C12H8O	000132-64-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId :
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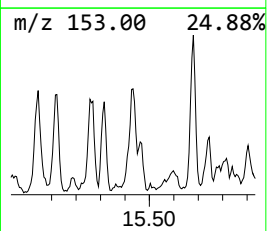
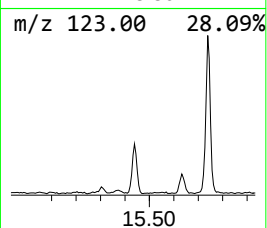
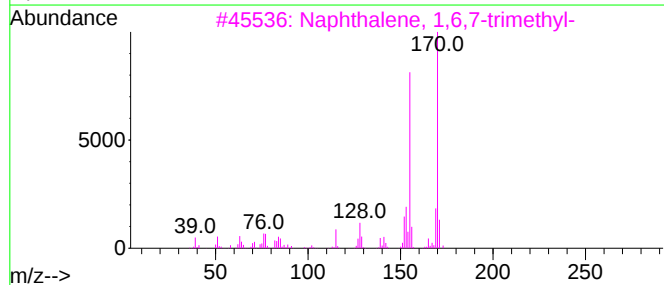
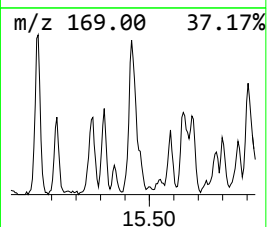
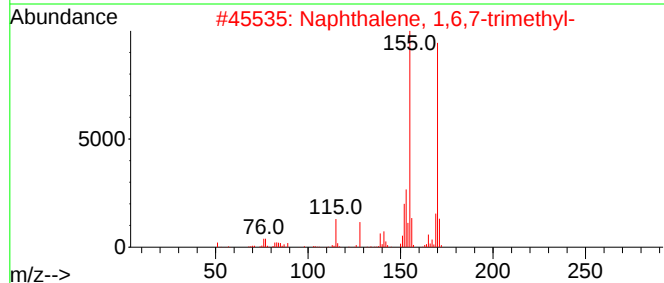
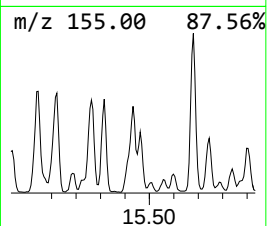
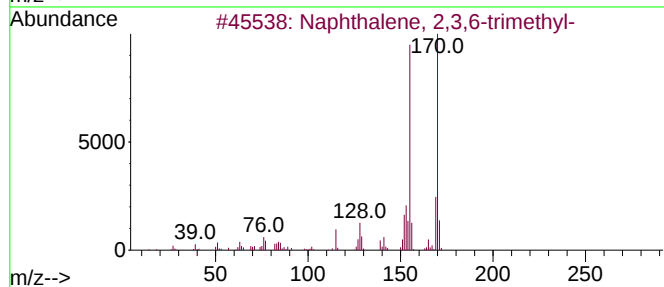
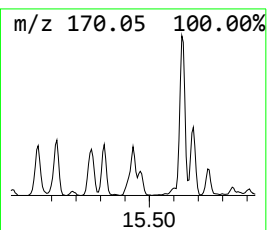
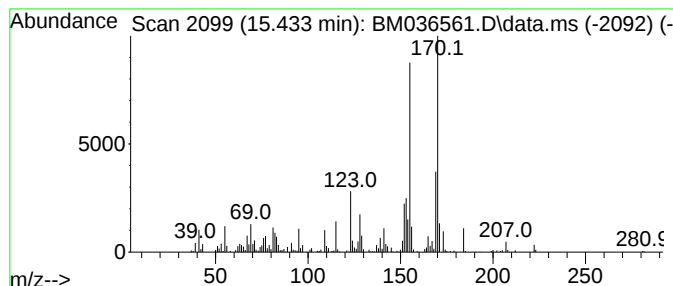
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1,6,7-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	2.19 ng/ul	490194	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
4			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	94
5			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Sample : N4601-12
Misc :
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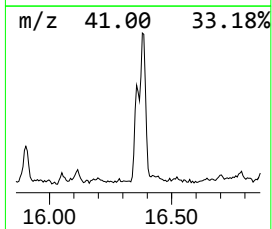
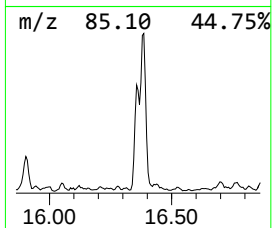
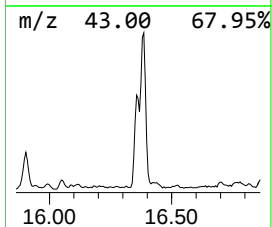
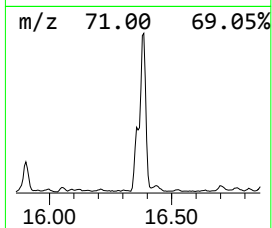
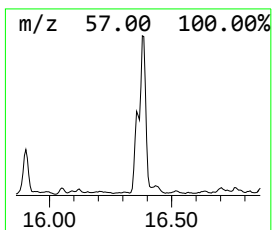
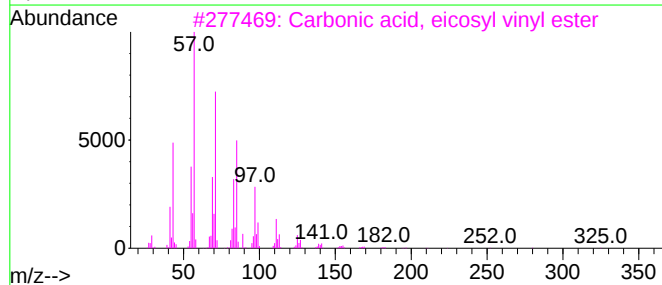
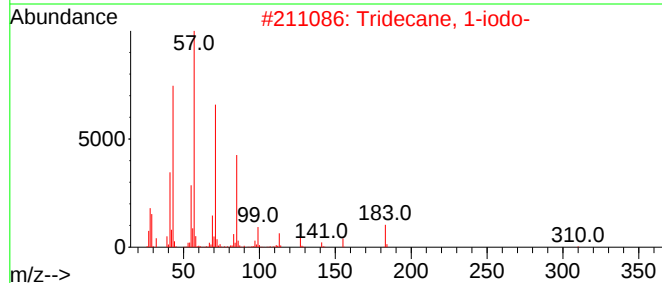
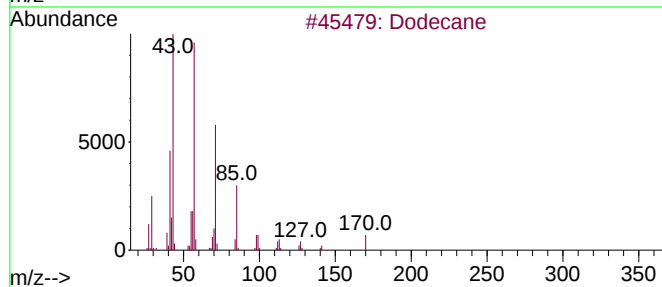
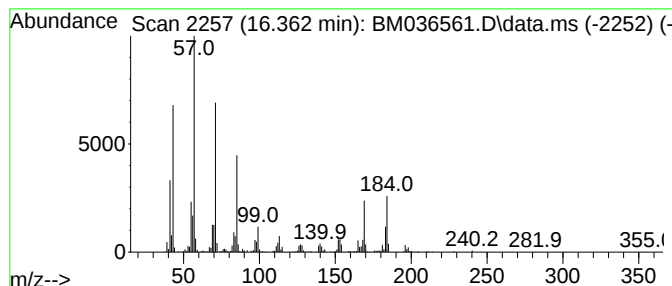
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.362	2.19 ng/ul	547599	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	64
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	58
3	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	58
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	53
5	Nonadecane	268	C19H40	000629-92-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId :
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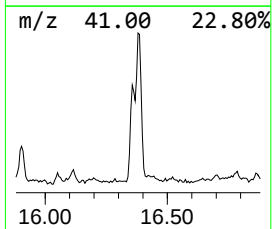
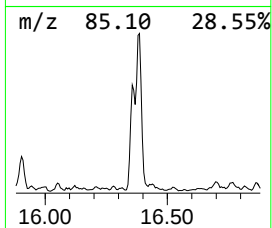
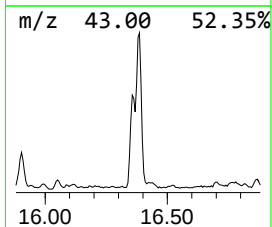
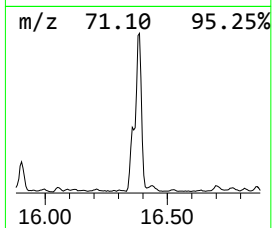
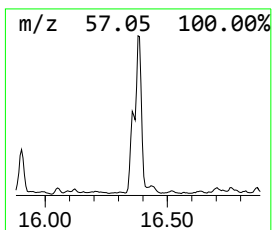
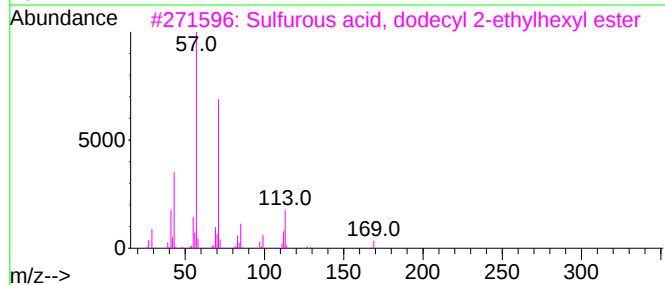
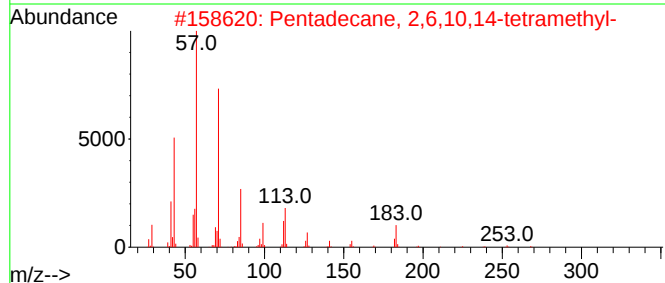
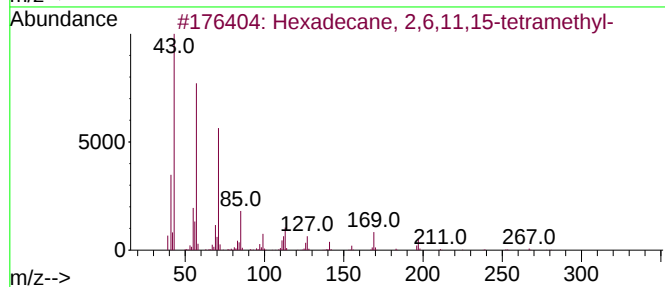
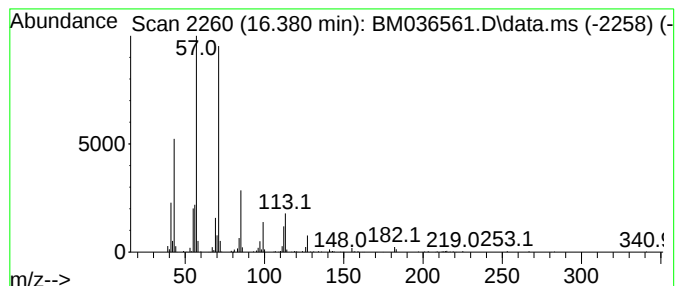
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.380	3.48 ng/ul	872082	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	83
2			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
3			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	72
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

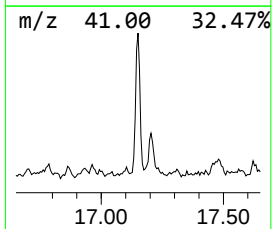
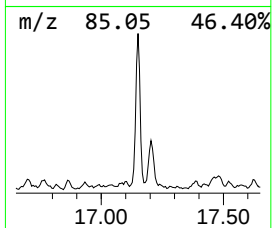
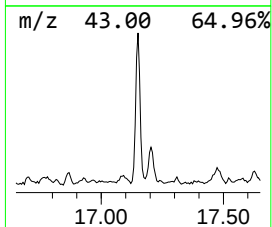
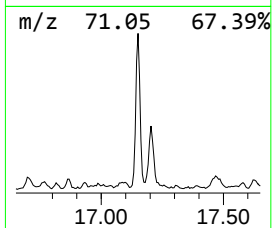
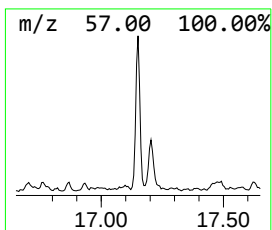
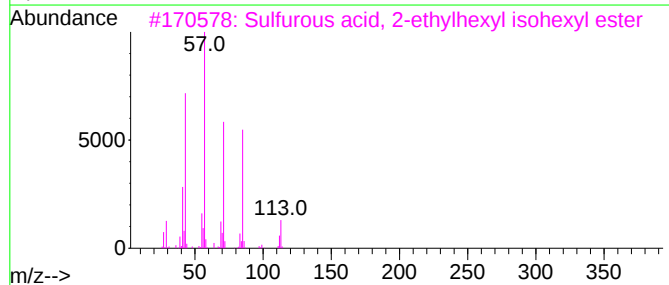
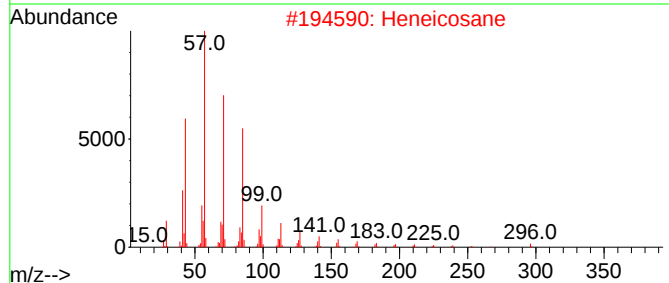
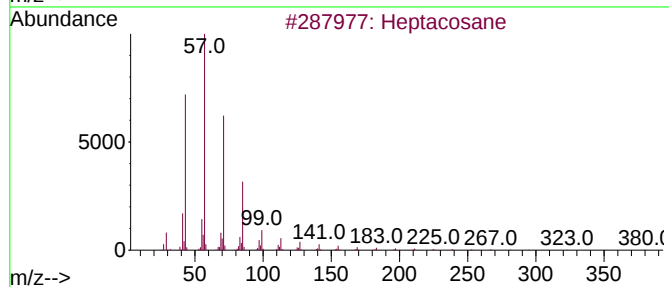
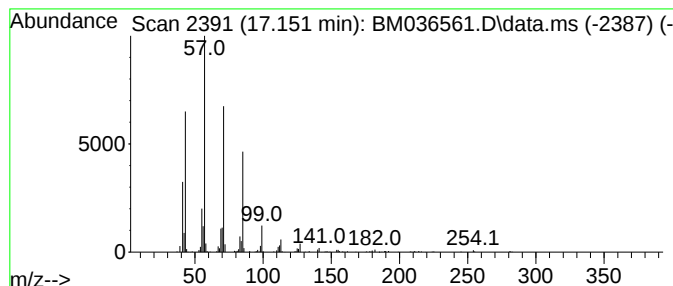
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.151	2.32 ng/ul	582121	Phenanthrene-d10		17.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Heneicosane		296	C21H44	000629-94-7	91
3	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	80
4	Tetracosane		338	C24H50	000646-31-1	72
5	Tetradecane		198	C14H30	000629-59-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

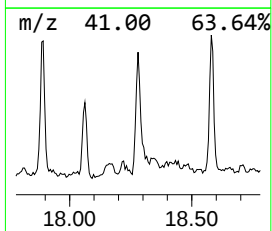
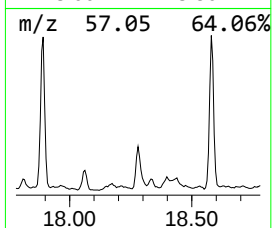
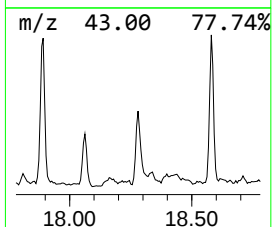
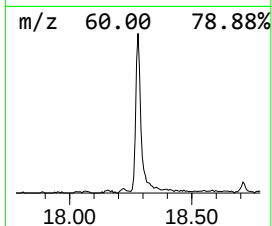
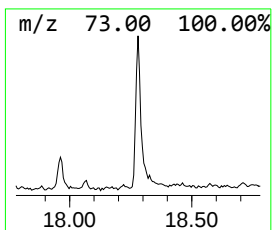
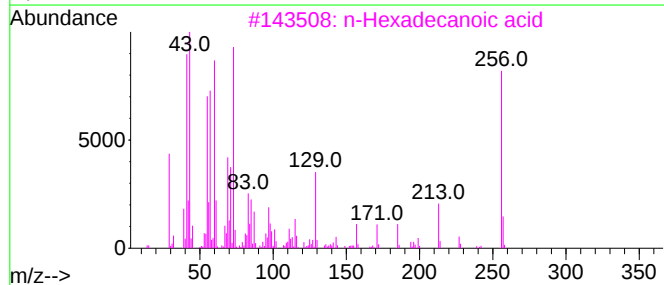
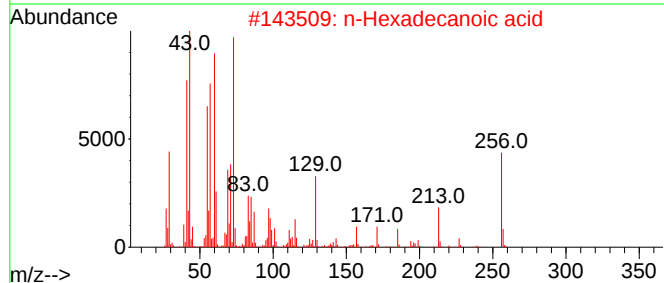
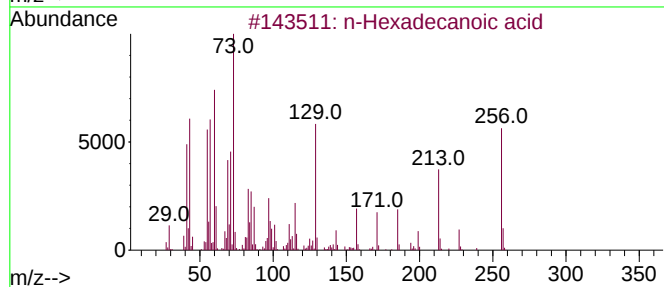
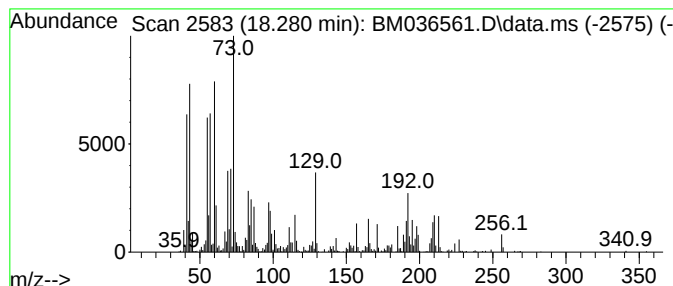
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.280	4.07 ng/ul	1019020	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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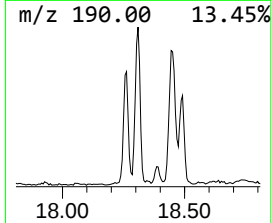
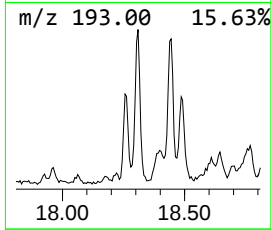
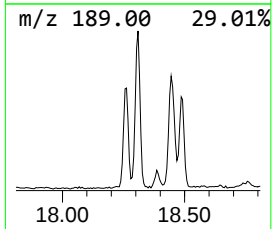
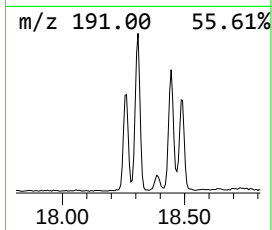
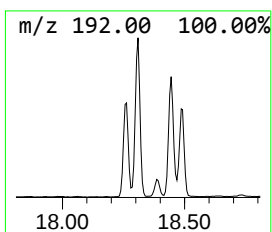
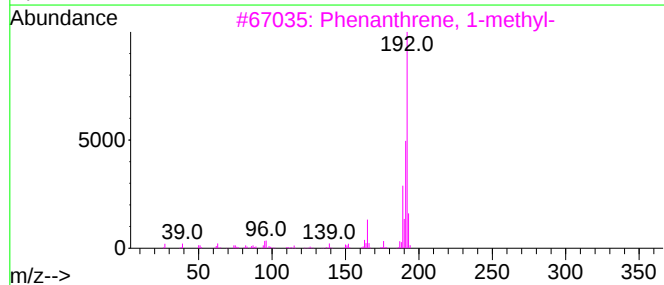
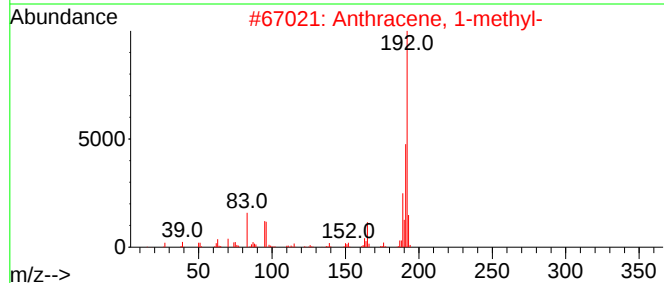
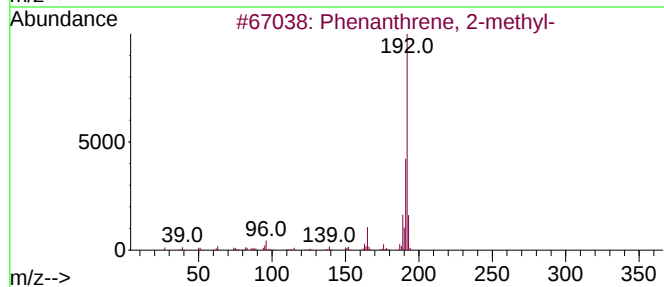
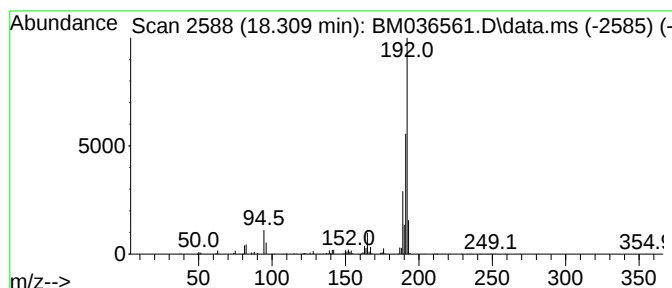
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.309	3.48 ng/ul	871290	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

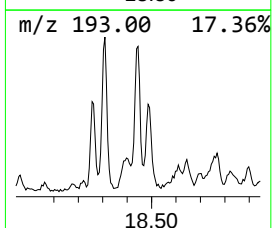
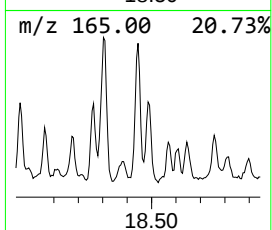
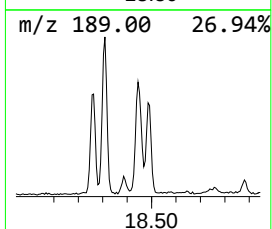
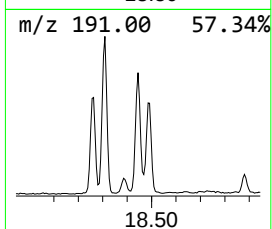
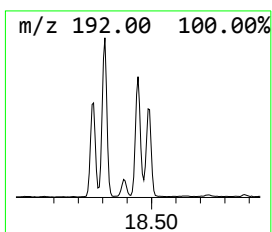
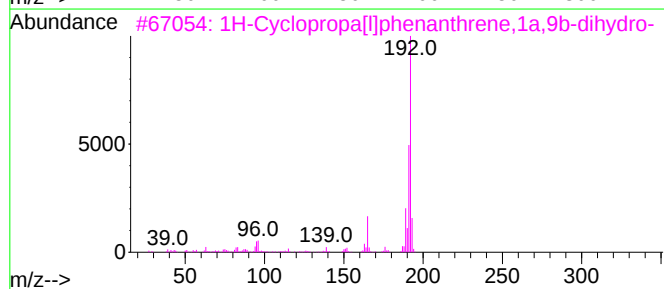
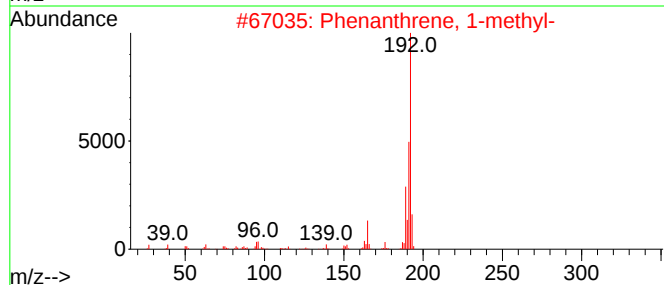
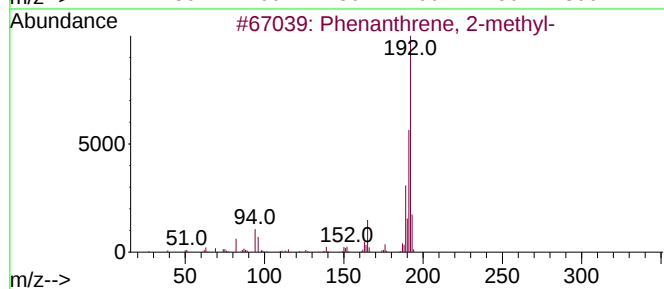
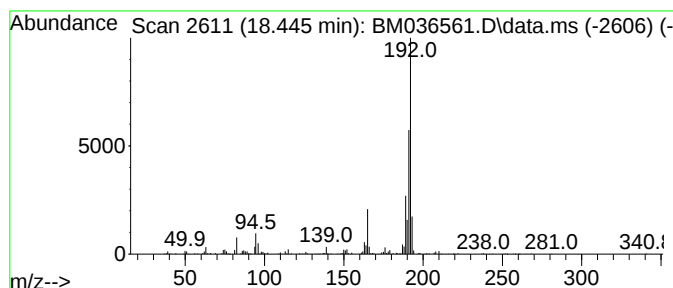
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 Phenanthrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.445	2.77 ng/ul	693842	Phenanthrene-d10		17.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-		192	C15H12	000832-69-9	96
3	1H-Cyclopropa[1]phenanthrene,1a,...		192	C15H12	000949-41-7	96
4	1H-Indene, 2-phenyl-		192	C15H12	004505-48-0	94
5	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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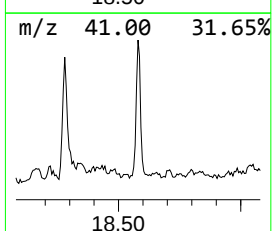
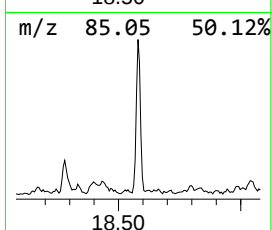
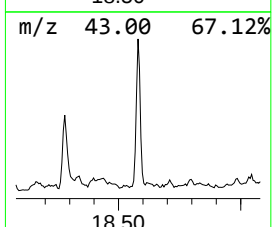
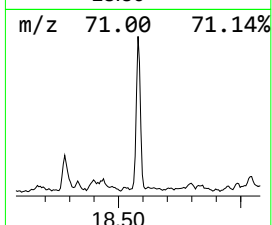
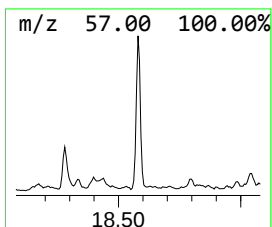
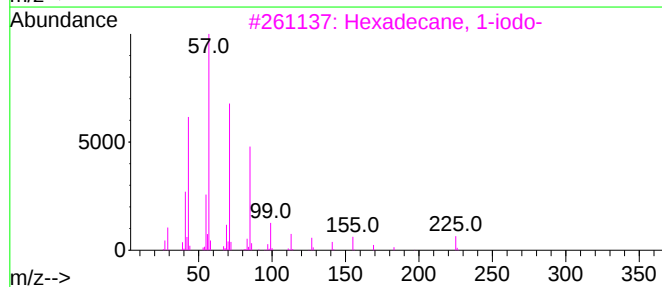
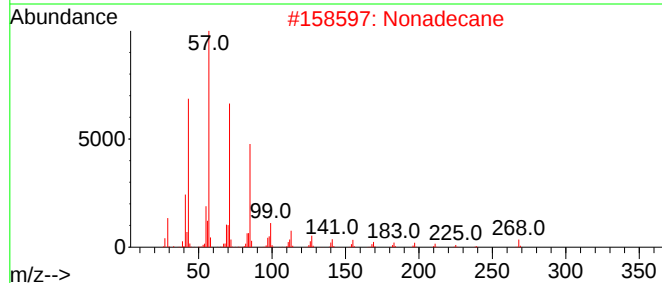
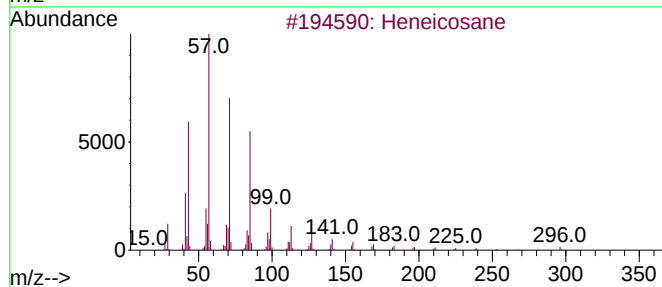
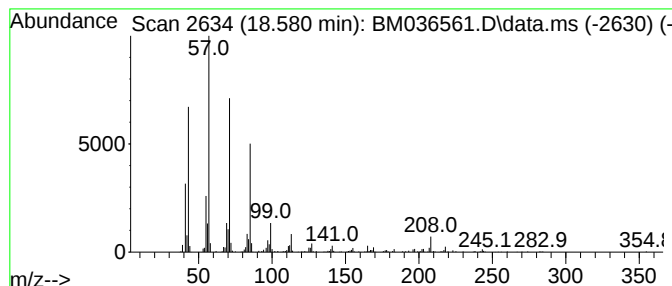
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.580	2.10 ng/ul	526198	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Nonadecane	268	C19H40	000629-92-5	91
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
4		Hexadecane	226	C16H34	000544-76-3	86
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

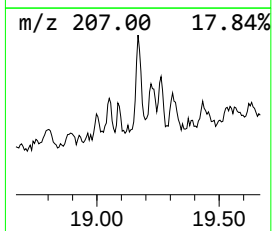
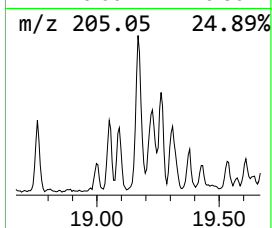
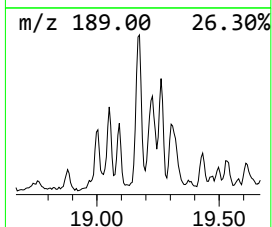
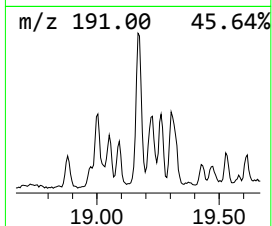
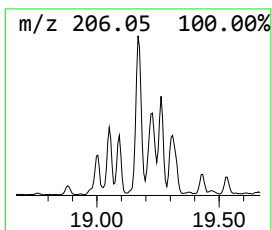
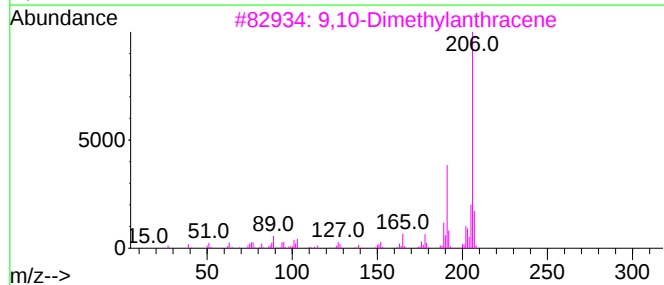
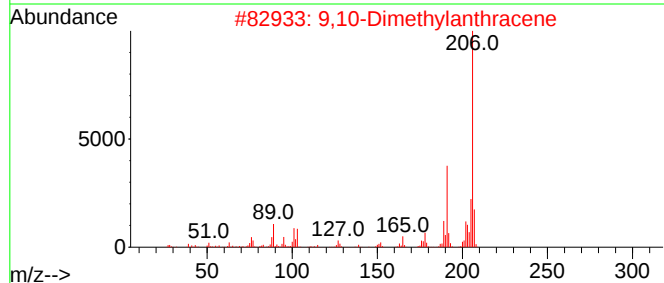
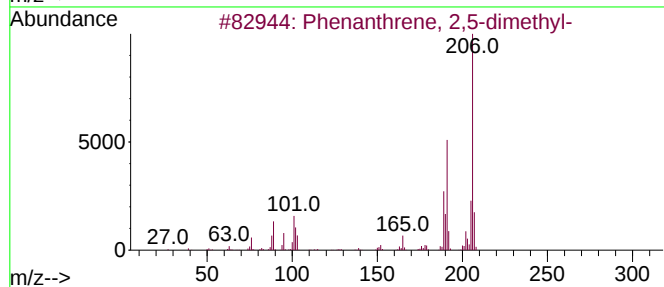
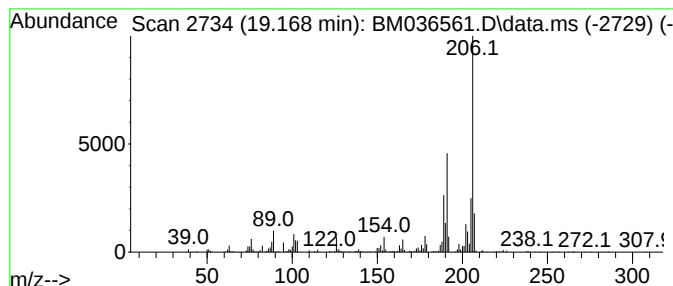
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 Phenanthrene, 2,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.168	3.09 ng/ul	773994	Phenanthrene-d10		17.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	93
2	9,10-Dimethylantracene		206	C16H14	000781-43-1	91
3	9,10-Dimethylantracene		206	C16H14	000781-43-1	91
4	Phenanthrene, 1,7-dimethyl-		206	C16H14	000483-87-4	90
5	9,10-Dimethylantracene		206	C16H14	000781-43-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

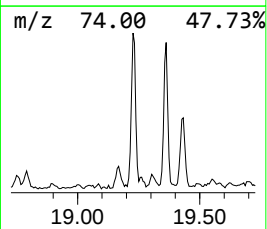
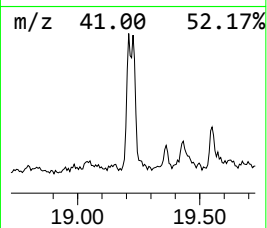
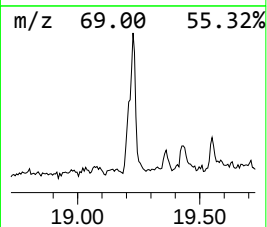
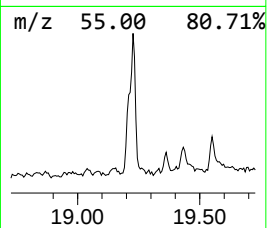
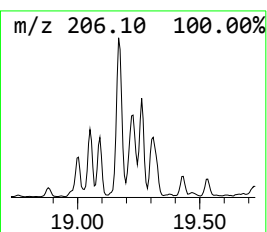
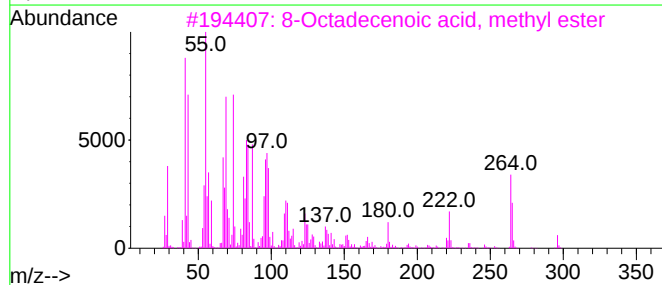
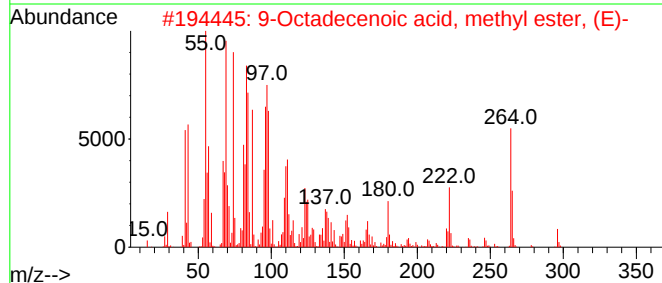
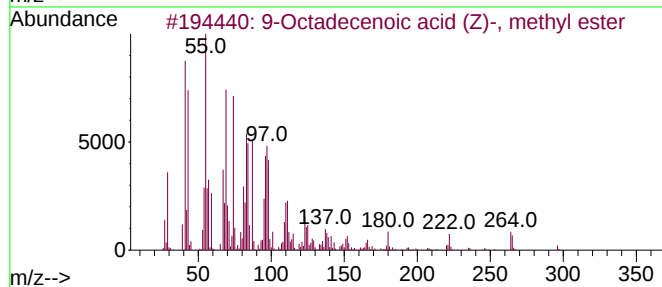
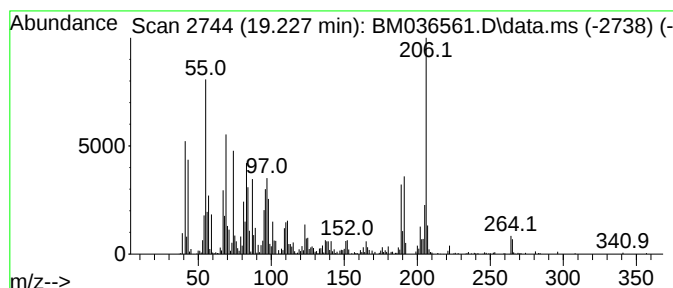
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 9-Octadecenoic acid (Z)-, m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.227	5.28 ng/ul	1322770	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	96
2	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	94
3	8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	87
4	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	87
5	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

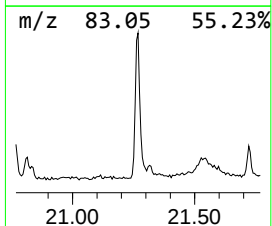
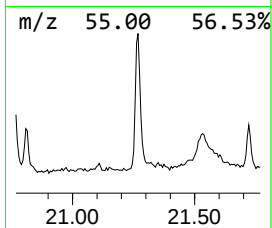
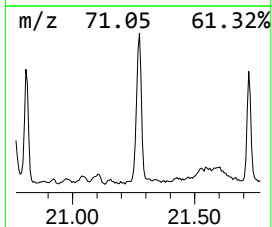
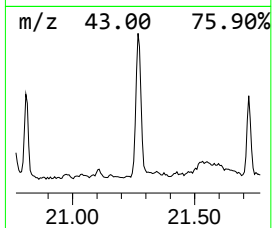
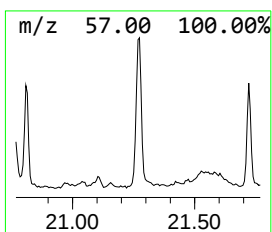
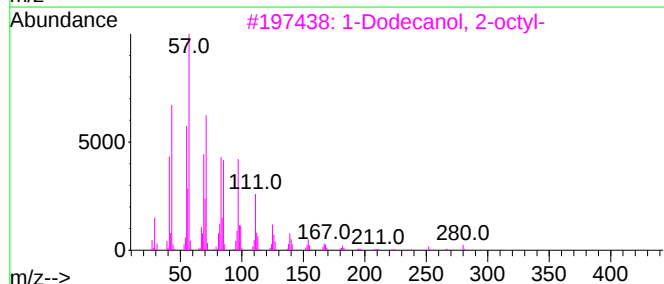
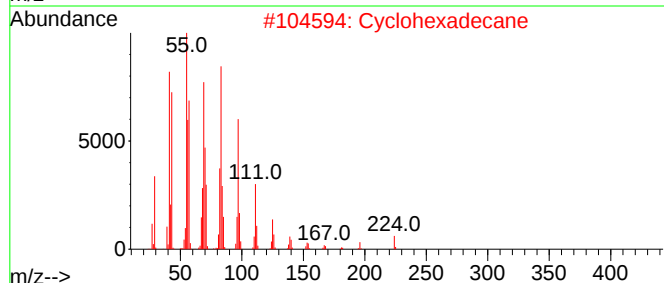
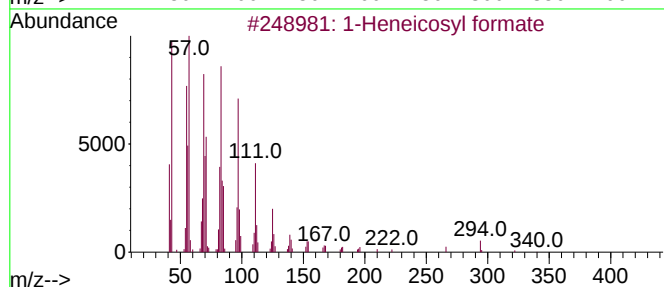
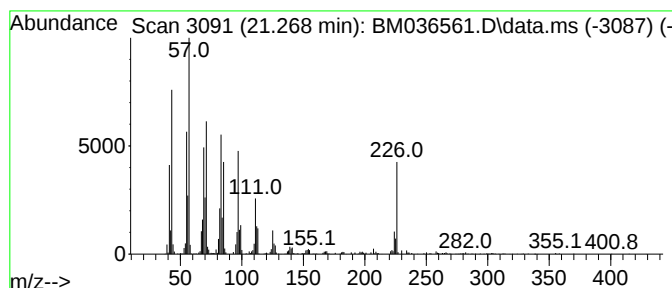
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 1-Heneicosyl formate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.268	3.49 ng/ul	1281450	Chrysene-d12		21.550	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate		340	C22H44O2	077899-03-7	89
2	Cyclohexadecane		224	C16H32	000295-65-8	89
3	1-Dodecanol, 2-octyl-		298	C20H42O	005333-42-6	81
4	Trihexadecyl borate		735	C48H99BO3	002665-11-4	76
5	Ethanol, 2-(tetradecyloxy)-		258	C16H34O2	002136-70-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

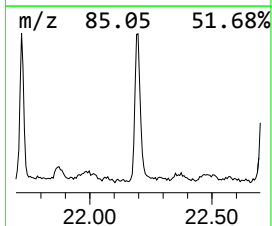
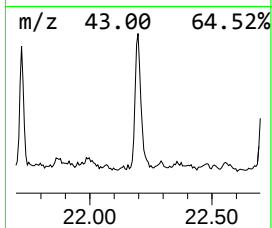
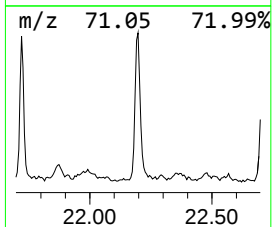
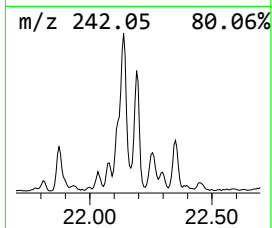
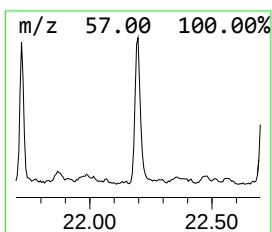
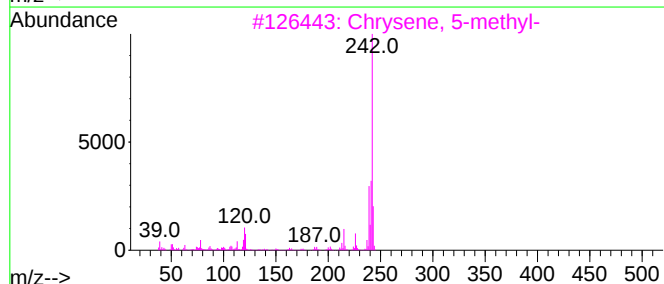
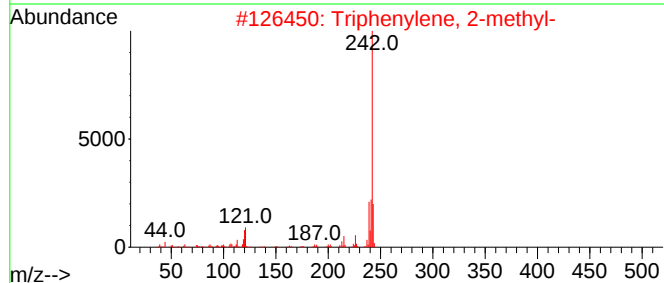
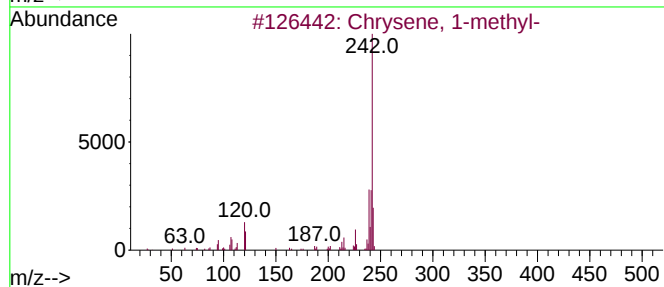
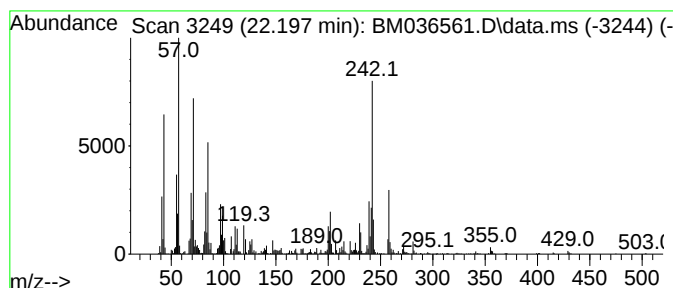
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 19 Chrysene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.197	3.80 ng/ul	1392950	Chrysene-d12		21.550	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Chrysene, 1-methyl-		242	C19H14	003351-28-8	92
2	Triphenylene, 2-methyl-		242	C19H14	001705-84-6	78
3	Chrysene, 5-methyl-		242	C19H14	003697-24-3	70
4	Hexadecane		226	C16H34	000544-76-3	66
5	Tetradecane, 1-chloro-		232	C14H29Cl	002425-54-9	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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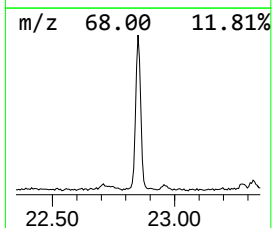
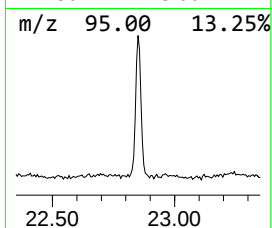
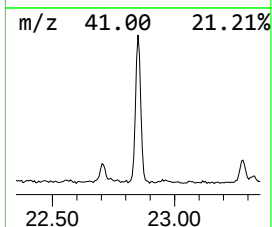
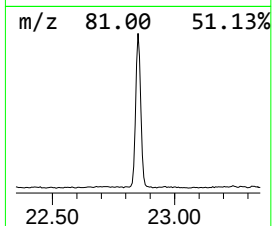
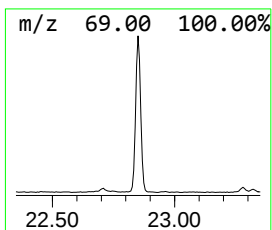
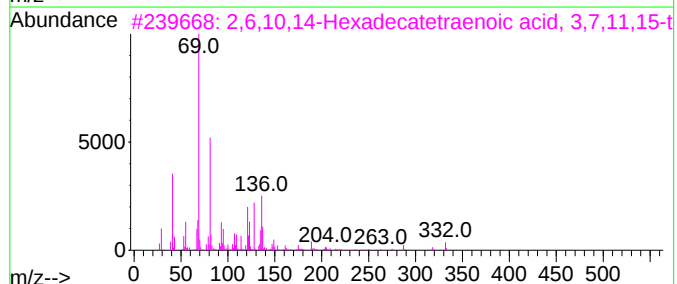
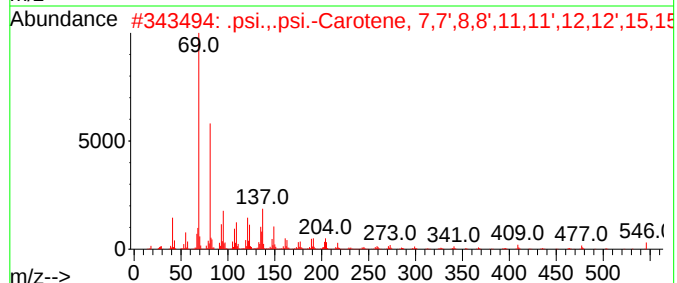
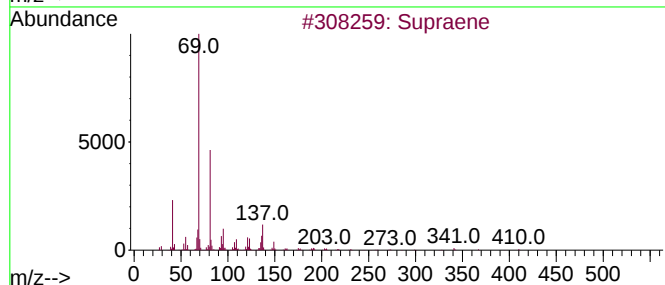
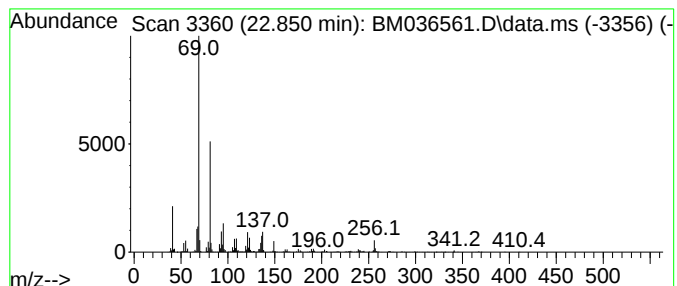
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Supraene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.850	9.43 ng/ul	2380510	Perylene-d12	23.997

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			.psi.,.psi.-Carotene, 7,7',8,8',...	547	C40H66	000502-62-5	83
3			2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	64
4			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	50
5			Methanone, dicyclopropyl-	110	C7H10O	001121-37-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
Operator : CG/JU
Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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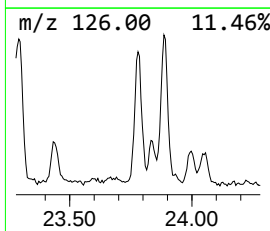
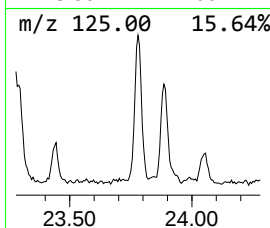
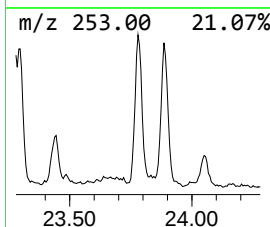
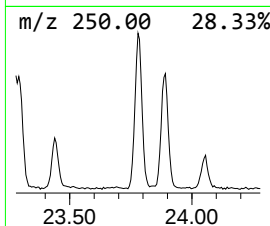
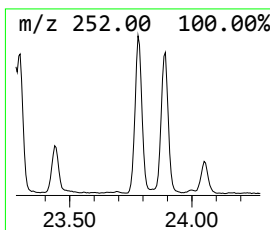
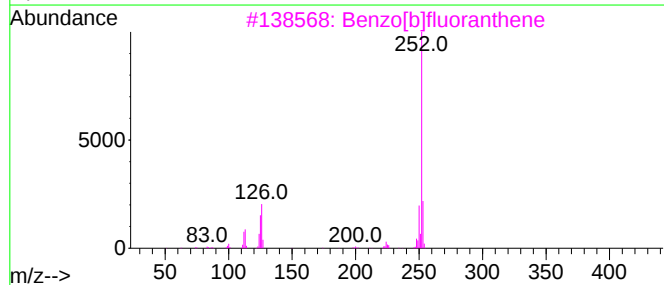
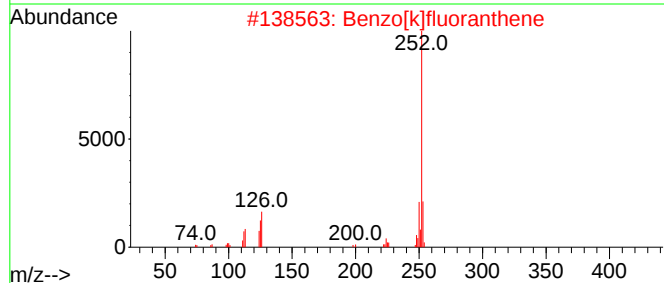
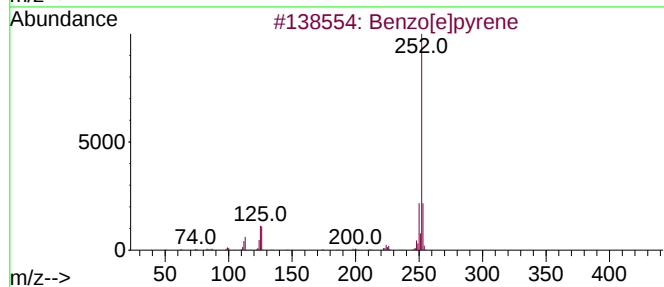
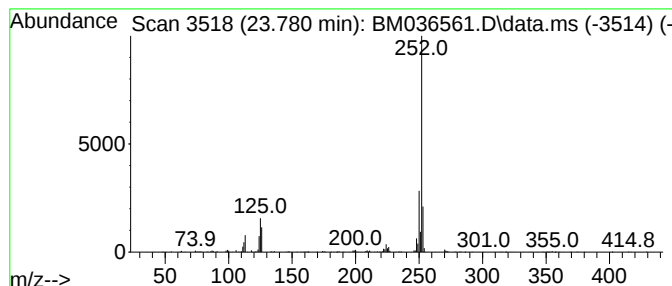
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzo[e]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.779	2.46 ng/ul	620566	Perylene-d12	23.997

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036561.D
Acq On : 16 Sep 2022 14:10
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Sample : N4601-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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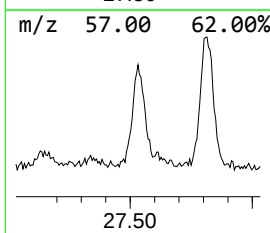
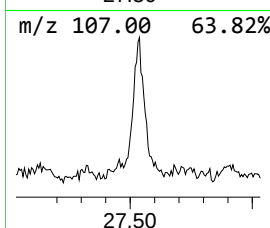
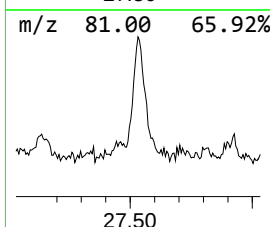
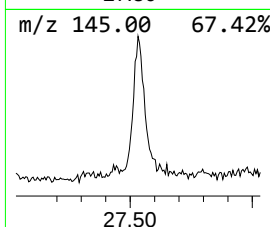
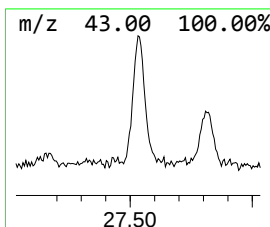
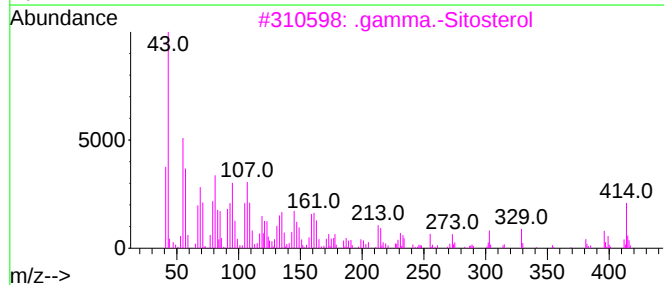
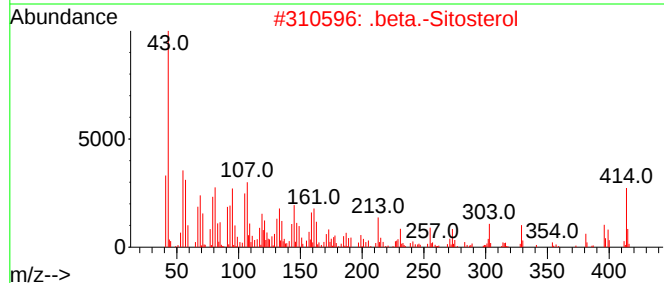
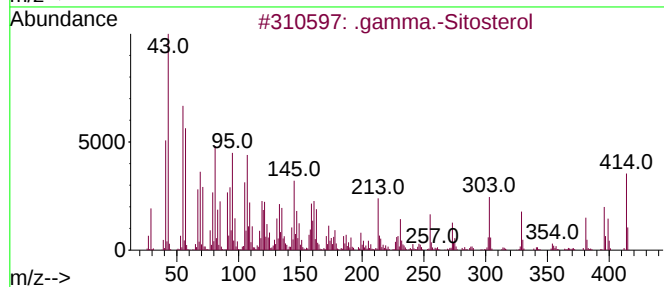
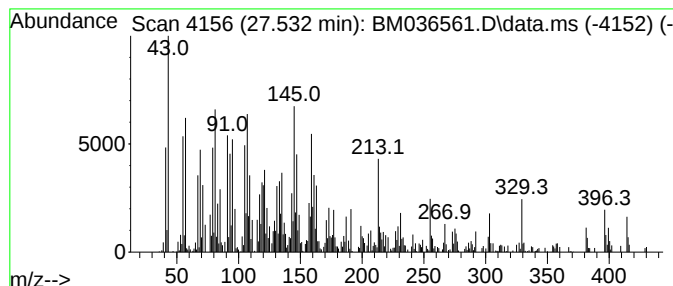
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 .gamma.-Sitosterol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.532	4.94 ng/ul	1247430	Perylene-d12	23.997

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	89	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	78	
3	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	64	
4	Stigmasta-3,5-diene	396	C29H48	004970-37-0	49		
5	Campesterol	400	C28H48O	000474-62-4	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036561.D
 Acq On : 16 Sep 2022 14:10
 Operator : CG/JU
 Sample : N4601-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.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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o-Xylene	6.022	2.2	ng/ul	224753	1	8.028	2005150	20.0
Ethanol, 2-(2-b...	10.610	7.7	ng/ul	1238600	2	10.839	3212620	20.0
(DEL) Alkane: S...	11.869	2.1	ng/ul	343280	2	10.839	3212620	20.0
(DEL) Alkane: S...	13.486	2.1	ng/ul	463808	3	14.645	4481470	20.0
Naphthalene, 2,...	13.974	2.6	ng/ul	586089	3	14.645	4481470	20.0
Naphthalene, 2,...	15.045	2.2	ng/ul	487731	3	14.645	4481470	20.0
Naphthalene, 1,...	15.433	2.2	ng/ul	490194	3	14.645	4481470	20.0
(DEL) Alkane: S...	16.362	2.2	ng/ul	547599	4	17.386	5009030	20.0
(DEL) Alkane: S...	16.380	3.5	ng/ul	872082	4	17.386	5009030	20.0
(DEL) Alkane: S...	17.151	2.3	ng/ul	582121	4	17.386	5009030	20.0
n-Hexadecanoic ...	18.280	4.1	ng/ul	1019020	4	17.386	5009030	20.0
Phenanthrene, 2...	18.309	3.5	ng/ul	871290	4	17.386	5009030	20.0
Phenanthrene, 1...	18.445	2.8	ng/ul	693842	4	17.386	5009030	20.0
(DEL) Alkane: S...	18.580	2.1	ng/ul	526198	4	17.386	5009030	20.0
Phenanthrene, 2...	19.168	3.1	ng/ul	773994	4	17.386	5009030	20.0
9-Octadecenoic ...	19.227	5.3	ng/ul	1322770	4	17.386	5009030	20.0
1-Heneicosyl fo...	21.268	3.5	ng/ul	1281450	5	21.550	7340390	20.0
Chrysene, 1-met...	22.197	3.8	ng/ul	1392950	5	21.550	7340390	20.0
Supraene	22.850	9.4	ng/ul	2380510	6	23.997	5049920	20.0
Benzo[e]pyrene	23.779	2.5	ng/ul	620566	6	23.997	5049920	20.0
.gamma.-Sitosterol	27.532	4.9	ng/ul	1247430	6	23.997	5049920	20.0