

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036562.D
 Acq On : 16 Sep 2022 14:46
 Operator : CG/JU
 Sample : N4601-14
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5755

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

Signal : TIC: BM036562.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	27	rVB	12195874	16466206	100.00%	11.586%
2	6.022	493	499	508	rVB3	86058	160970	0.98%	0.113%
3	7.169	688	694	703	rVV	1180629	1986298	12.06%	1.398%
4	7.351	719	725	732	rVV	1040178	1582911	9.61%	1.114%
5	7.551	749	759	768	rBV	1169830	1839848	11.17%	1.295%
6	7.669	773	779	789	rVV2	76693	192643	1.17%	0.136%
7	8.028	832	840	846	rBV	1267584	2088107	12.68%	1.469%
8	8.728	952	959	967	rVB	1177440	1906643	11.58%	1.342%
9	9.186	1031	1037	1043	rVV	1140904	1844584	11.20%	1.298%
10	9.263	1047	1050	1061	rVB2	92047	185939	1.13%	0.131%
11	9.916	1154	1161	1167	rBV	863249	1432645	8.70%	1.008%
12	10.451	1243	1252	1261	rVB	1579038	2592328	15.74%	1.824%
13	10.610	1271	1279	1288	rBV	658077	1007654	6.12%	0.709%
14	10.839	1309	1318	1323	rBV	1910731	3326939	20.20%	2.341%
15	10.886	1323	1326	1334	rVB	219460	338778	2.06%	0.238%
16	10.969	1334	1340	1345	rBV	754680	1316552	8.00%	0.926%
17	11.869	1488	1493	1501	rBV	243642	378736	2.30%	0.266%
18	12.257	1554	1559	1571	rBV	179858	306237	1.86%	0.215%
19	12.492	1593	1599	1605	rVB	413844	688050	4.18%	0.484%
20	12.704	1630	1635	1644	rVV	383862	607917	3.69%	0.428%
21	13.204	1716	1720	1725	rVB	183514	249292	1.51%	0.175%
22	13.486	1762	1768	1776	rBV	353051	546582	3.32%	0.385%
23	13.816	1819	1824	1826	rBV	145459	228664	1.39%	0.161%
24	13.974	1846	1851	1856	rVV	465259	763638	4.64%	0.537%
25	14.027	1856	1860	1861	rVV	351168	420108	2.55%	0.296%
26	14.063	1861	1866	1873	rVV	3047617	4478209	27.20%	3.151%
27	14.139	1875	1879	1883	rVV	332519	461823	2.80%	0.325%
28	14.210	1883	1891	1894	rVV2	279682	486013	2.95%	0.342%
29	14.245	1894	1897	1902	rVV2	107838	168758	1.02%	0.119%
30	14.345	1908	1914	1929	rVB	3707776	5449716	33.10%	3.835%
31	14.551	1945	1949	1955	rVB	305878	406709	2.47%	0.286%
32	14.651	1956	1966	1972	rBV	3009621	4661564	28.31%	3.280%
33	14.821	1990	1995	2000	rBV	1149985	1566254	9.51%	1.102%
34	15.045	2028	2033	2040	rVB	383202	636861	3.87%	0.448%
35	15.121	2040	2046	2050	rVB	203449	303033	1.84%	0.213%
36	15.174	2050	2055	2062	rVB5	88904	186938	1.14%	0.132%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
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37	15.263	2063	2070	2075	rBV	210716	371407	2.26%	0.261%
38	15.315	2075	2079	2083	rVB	190998	245043	1.49%	0.172%
39	15.433	2093	2099	2103	rBV3	264726	484803	2.94%	0.341%
40	15.498	2107	2110	2115	rVB	332996	409744	2.49%	0.288%
41	15.639	2126	2134	2139	rVV	4666054	6740984	40.94%	4.743%
42	15.680	2139	2141	2146	rVB	265605	314930	1.91%	0.222%
43	15.745	2146	2152	2157	rBV	1040045	1453526	8.83%	1.023%
44	15.904	2175	2179	2184	rVB	264756	384188	2.33%	0.270%
45	15.986	2185	2193	2202	rVB	282840	534419	3.25%	0.376%
46	16.133	2212	2218	2225	rVB3	238476	573091	3.48%	0.403%
47	16.221	2225	2233	2241	rVB4	142671	330260	2.01%	0.232%
48	16.362	2252	2257	2258	rBV2	519806	634050	3.85%	0.446%
49	16.386	2258	2261	2269	rVB	869535	1163128	7.06%	0.818%
50	16.921	2348	2352	2356	rBV3	217276	365696	2.22%	0.257%
51	17.039	2368	2372	2379	rVV5	166112	304757	1.85%	0.214%
52	17.115	2379	2385	2388	rVV	165097	341296	2.07%	0.240%
53	17.151	2388	2391	2397	rVV	458938	655837	3.98%	0.461%
54	17.204	2397	2400	2410	rVB4	238576	427857	2.60%	0.301%
55	17.386	2425	2431	2435	rBV	3671244	5166511	31.38%	3.635%
56	17.427	2435	2438	2442	rVV	931402	1224096	7.43%	0.861%
57	17.486	2442	2448	2457	rVB	5067513	7224253	43.87%	5.083%
58	17.574	2458	2463	2468	rBV3	113879	216366	1.31%	0.152%
59	17.698	2480	2484	2492	rVB3	221931	335164	2.04%	0.236%
60	17.892	2513	2517	2522	rVB	453010	564930	3.43%	0.398%
61	17.962	2526	2529	2533	rVB	164432	199454	1.21%	0.140%
62	18.062	2543	2546	2550	rVB	200708	217131	1.32%	0.153%
63	18.262	2576	2580	2582	rVV	420278	640574	3.89%	0.451%
64	18.309	2585	2588	2595	rVV	802086	1122322	6.82%	0.790%
65	18.386	2598	2601	2605	rVV3	107587	189787	1.15%	0.134%
66	18.445	2606	2611	2615	rVV	630995	963228	5.85%	0.678%
67	18.492	2615	2619	2625	rVB	468814	622405	3.78%	0.438%
68	18.580	2630	2634	2643	rVV	460041	660036	4.01%	0.464%
69	18.756	2660	2664	2668	rBV	341986	473591	2.88%	0.333%
70	18.792	2668	2670	2677	rVB4	121671	178096	1.08%	0.125%
71	19.056	2711	2715	2718	rBV2	163191	253192	1.54%	0.178%
72	19.174	2730	2735	2738	rBV	543636	815574	4.95%	0.574%
73	19.209	2738	2741	2747	rVV2	586016	975654	5.93%	0.687%
74	19.262	2747	2750	2754	rVV	293784	373741	2.27%	0.263%
75	19.309	2754	2758	2766	rVB2	268737	571925	3.47%	0.402%
76	19.433	2775	2779	2787	rVB	1229513	1584150	9.62%	1.115%

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77	19.498	2787	2790	2793	rVB	147538	159945	0.97%	0.113%
78	19.550	2793	2799	2802	rBV5	189562	328724	2.00%	0.231%
79	19.621	2808	2811	2817	rVB4	118503	204443	1.24%	0.144%
80	19.768	2830	2836	2849	rVV2	6299023	10623607	64.52%	7.475%
81	19.868	2849	2853	2859	rVB3	332016	562470	3.42%	0.396%
82	20.109	2890	2894	2899	rBV3	229579	521431	3.17%	0.367%
83	20.209	2907	2911	2914	rBV3	142969	233624	1.42%	0.164%
84	20.286	2921	2924	2927	rVB	180298	206068	1.25%	0.145%
85	20.321	2927	2930	2933	rVB	464026	503223	3.06%	0.354%
86	20.439	2946	2950	2955	rVB3	233290	433065	2.63%	0.305%
87	20.580	2972	2974	2978	rBV	124436	151110	0.92%	0.106%
88	20.768	3003	3006	3009	rBV	226061	228201	1.39%	0.161%
89	20.809	3010	3013	3021	rBV	430897	536234	3.26%	0.377%
90	21.027	3047	3050	3056	rVB	474623	710968	4.32%	0.500%
91	21.274	3087	3092	3097	rVB2	789997	1115173	6.77%	0.785%
92	21.550	3133	3139	3143	rBV	4730605	6721308	40.82%	4.729%
93	21.592	3143	3146	3155	rVB	627405	995891	6.05%	0.701%
94	21.721	3165	3168	3172	rBV	540302	687297	4.17%	0.484%
95	22.197	3245	3249	3255	rVB3	696220	970472	5.89%	0.683%
96	22.709	3332	3336	3343	rVB	394978	600103	3.64%	0.422%
97	23.250	3424	3428	3429	rBV	352920	449404	2.73%	0.316%
98	23.844	3522	3529	3535	rBV2	4026328	7707388	46.81%	5.423%
99	23.997	3548	3555	3563	rBV2	2699737	5344649	32.46%	3.761%
100	24.680	3667	3671	3679	rVB	279216	553376	3.36%	0.389%

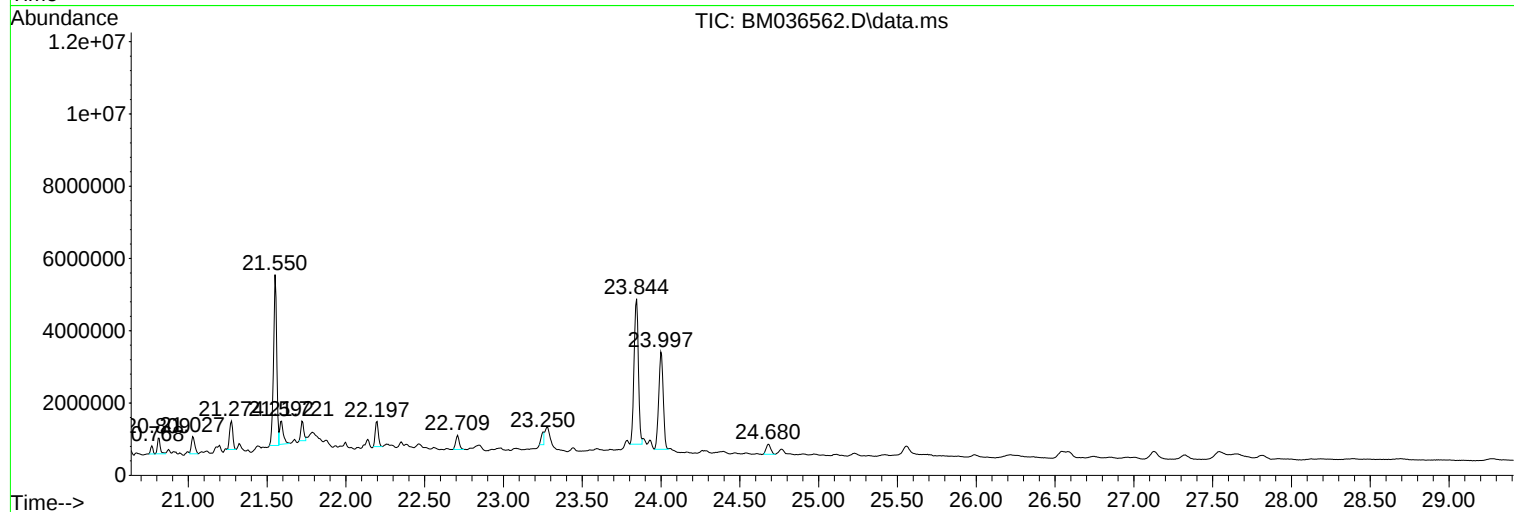
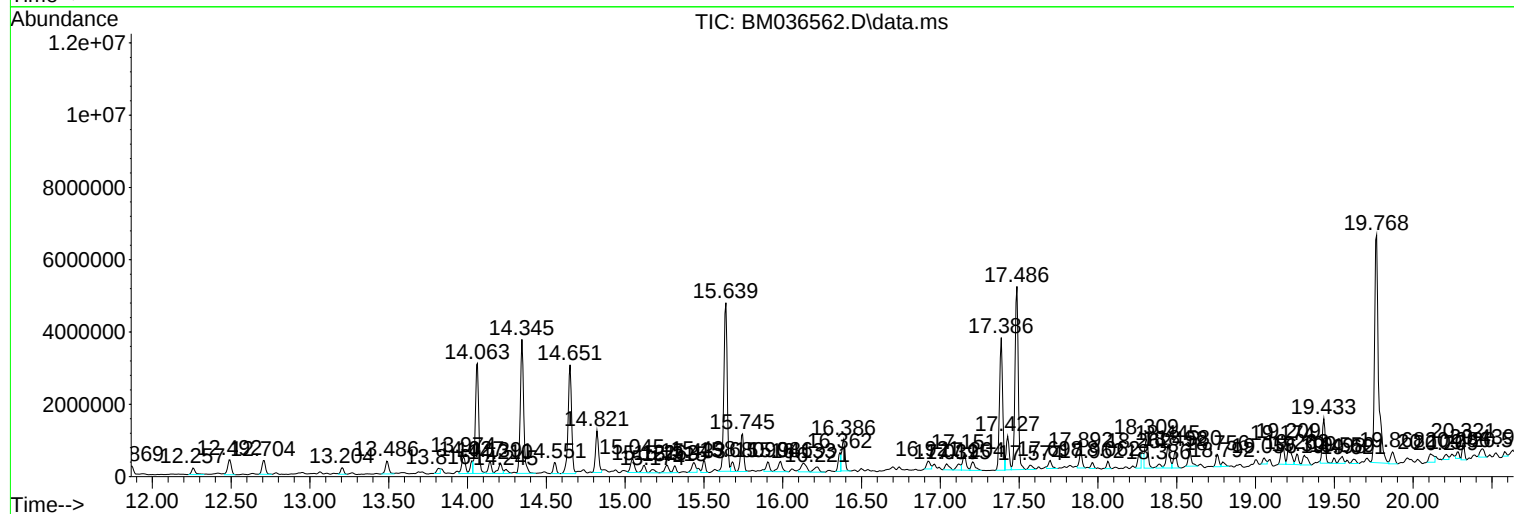
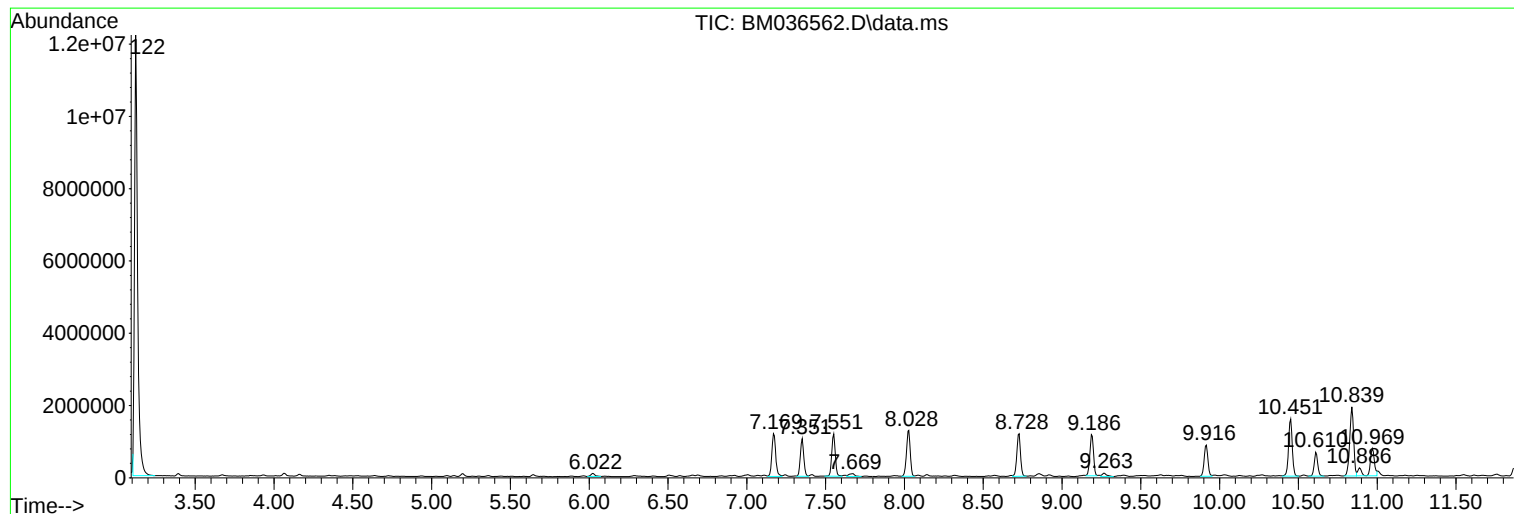
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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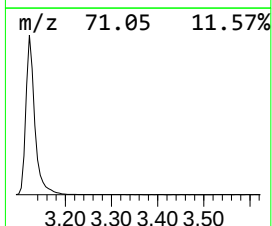
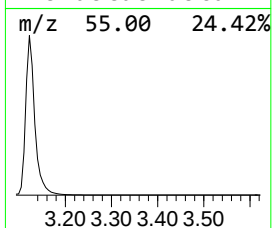
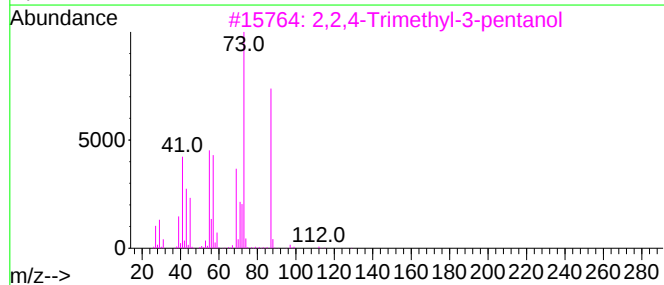
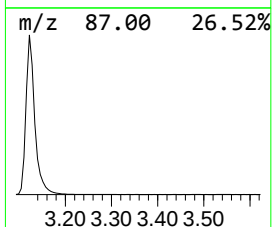
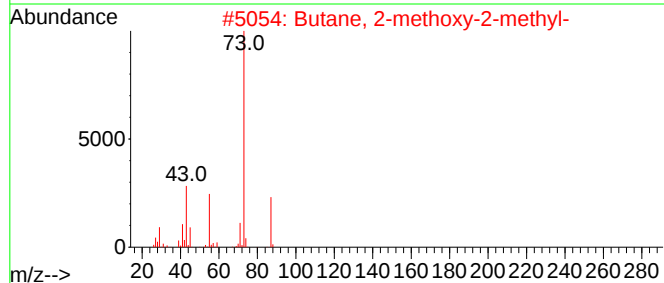
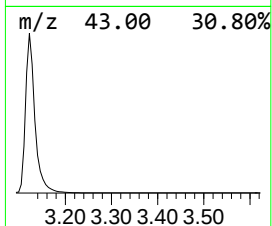
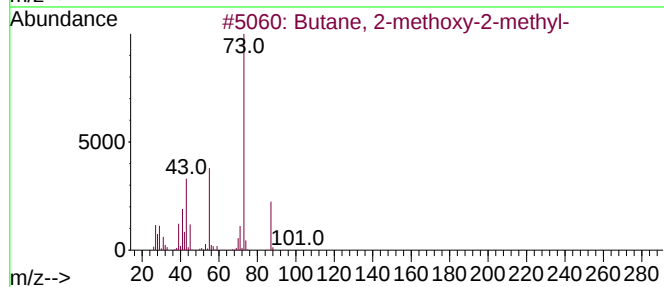
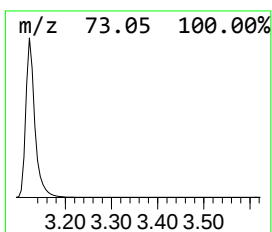
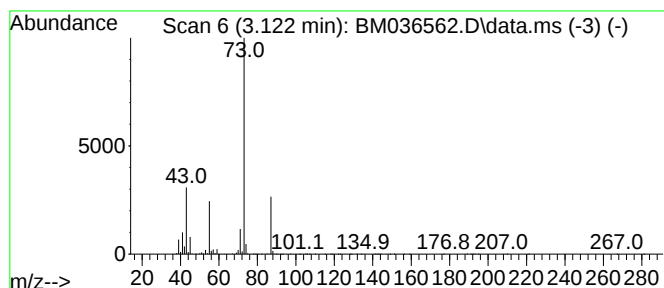
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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	157.71 ng/ul	16466200	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	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9		



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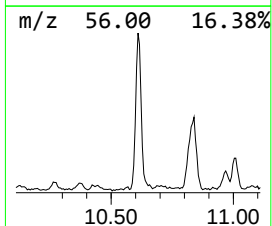
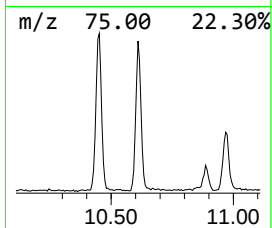
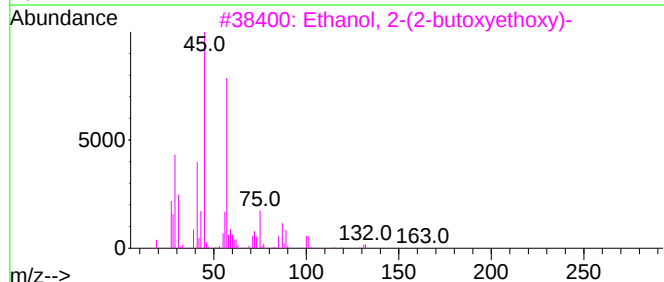
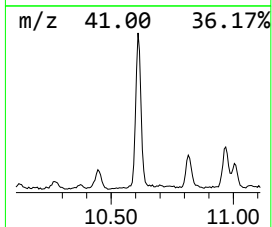
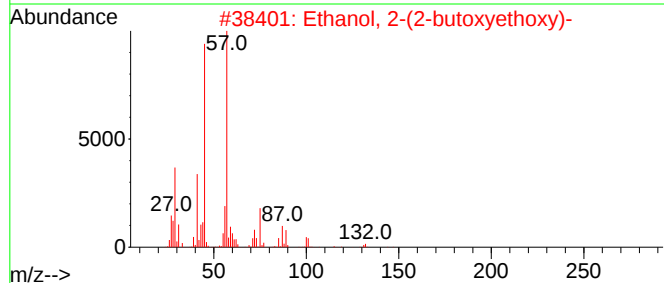
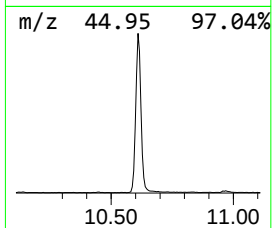
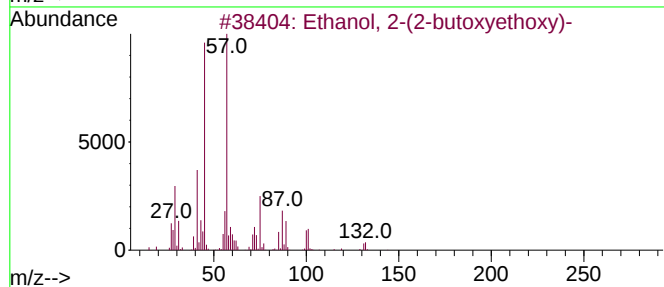
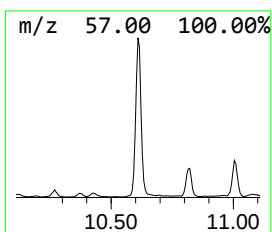
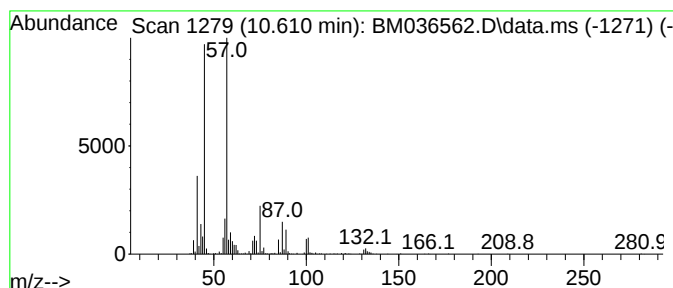
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.610	6.06 ng/ul	1007650	Naphthalene-d8	10.839	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78



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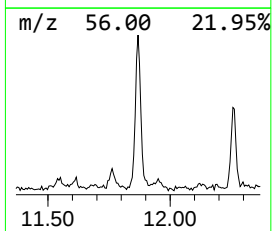
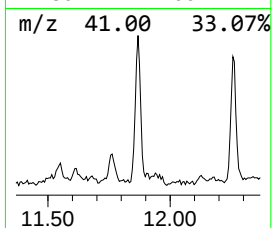
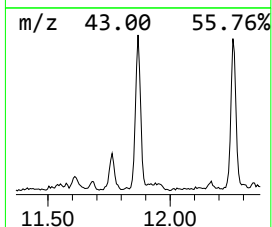
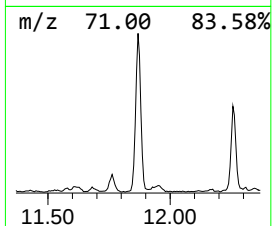
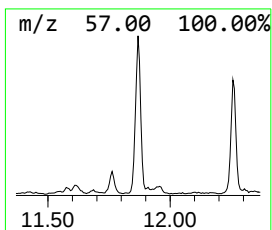
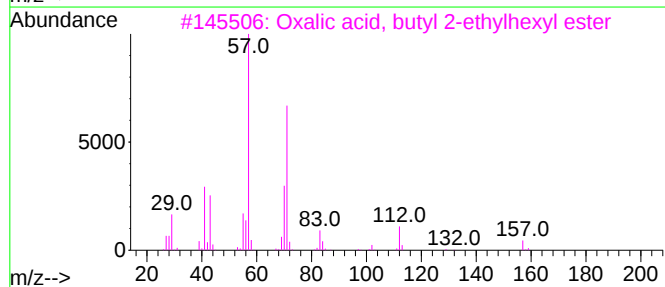
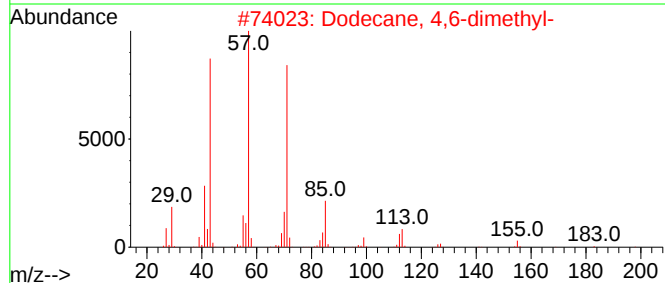
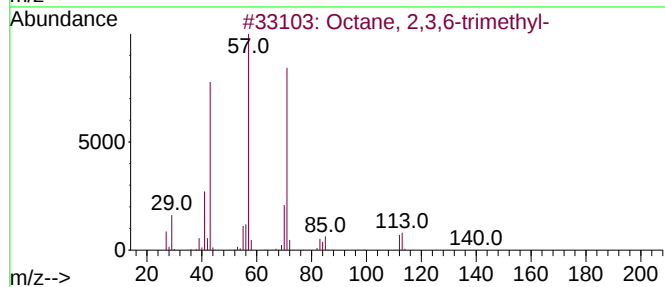
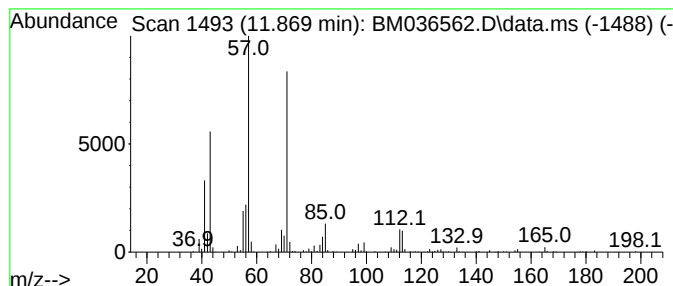
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.869	2.28 ng/ul	378736	Naphthalene-d8	10.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,3,6-trimethyl-			156	C11H24	062016-33-5	72
2	Dodecane, 4,6-dimethyl-			198	C14H30	061141-72-8	72
3	Oxalic acid, butyl 2-ethylhexyl ...			258	C14H26O4	1000309-38-6	59
4	Oxalic acid, bis(2-ethylhexyl) e...			314	C18H34O4	1000309-39-0	59
5	Oxalic acid, 2-ethylhexyl isobut...			258	C14H26O4	1010309-38-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

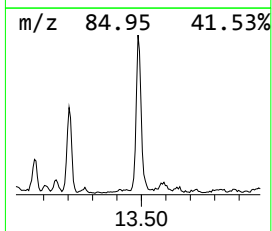
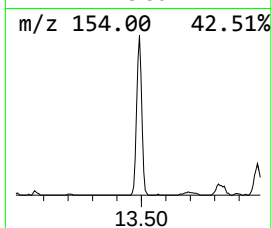
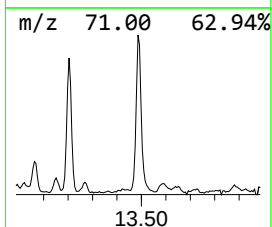
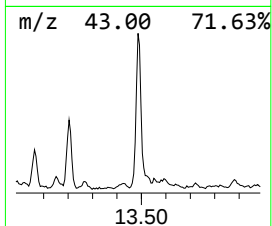
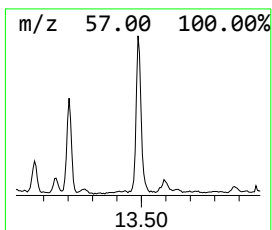
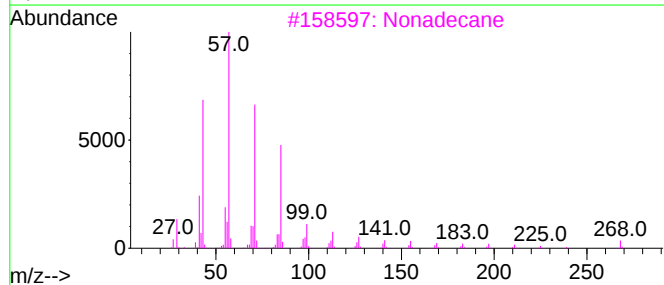
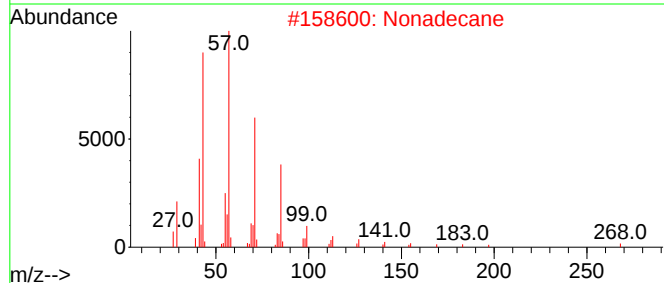
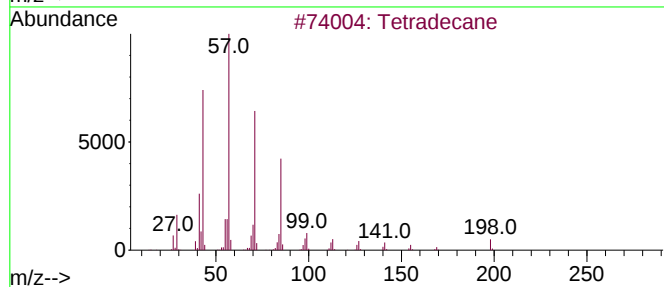
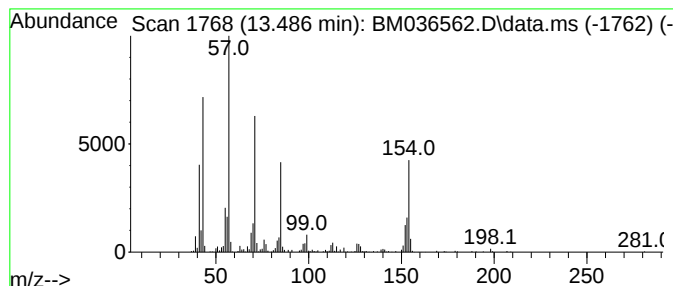
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.486	2.35 ng/ul	546582	Acenaphthene-d10		14.651	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	90
2	Nonadecane		268	C19H40	000629-92-5	53
3	Nonadecane		268	C19H40	000629-92-5	49
4	Tridecane		184	C13H28	000629-50-5	49
5	Tetradecane		198	C14H30	000629-59-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
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Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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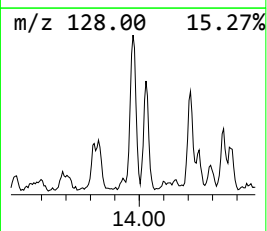
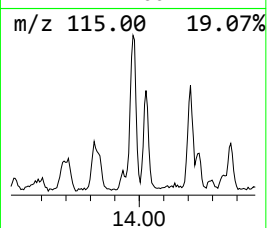
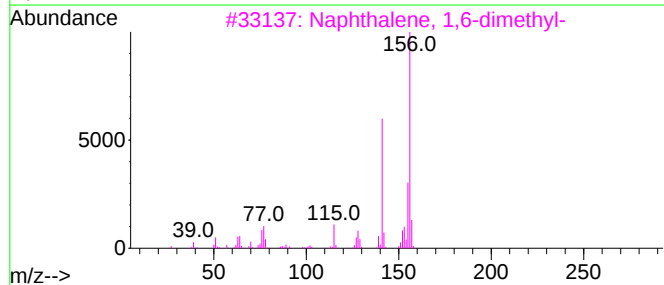
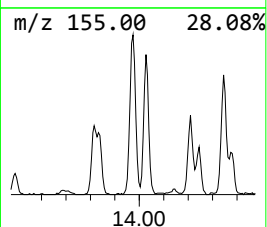
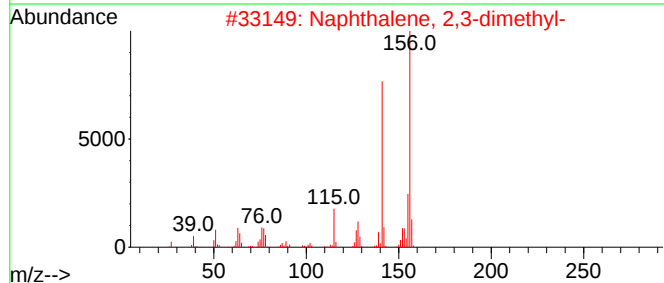
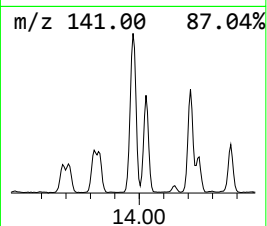
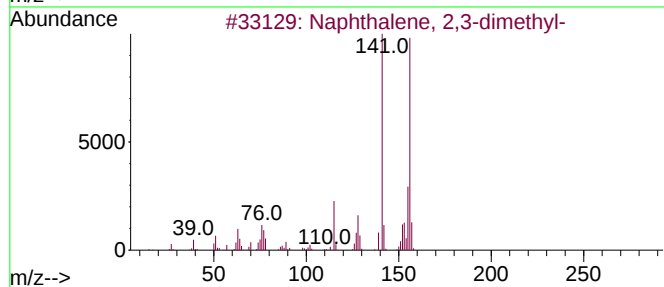
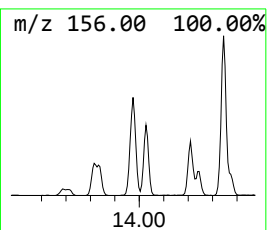
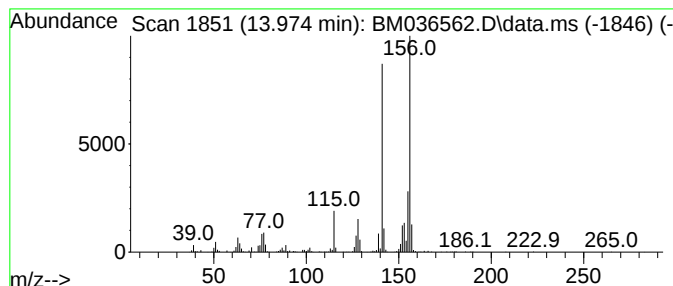
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.974	3.28 ng/ul	763638	Acenaphthene-d10	14.651

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, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

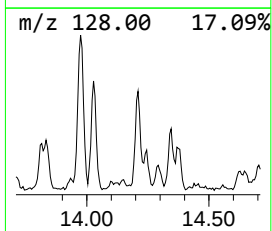
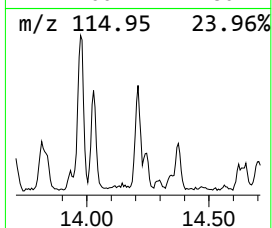
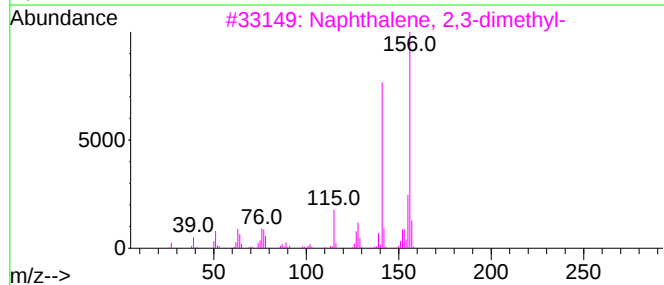
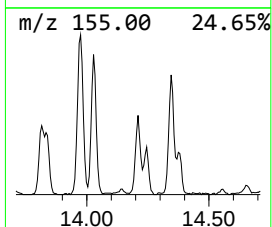
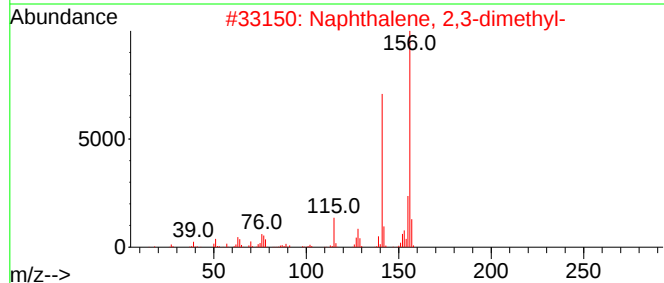
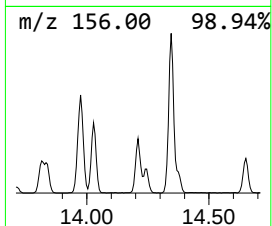
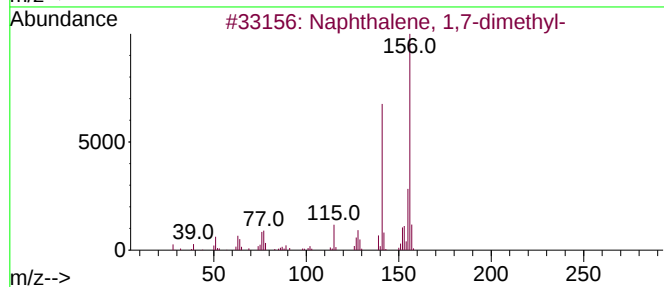
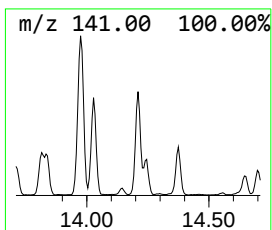
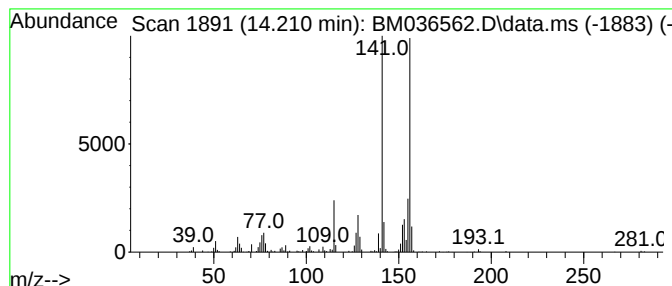
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 1,7-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.210	2.09 ng/ul	486013	Acenaphthene-d10	14.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
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Sample : N4601-14
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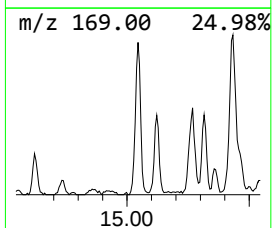
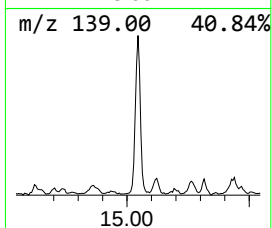
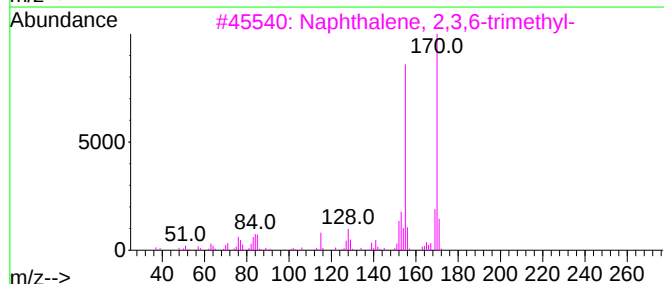
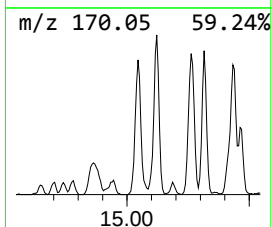
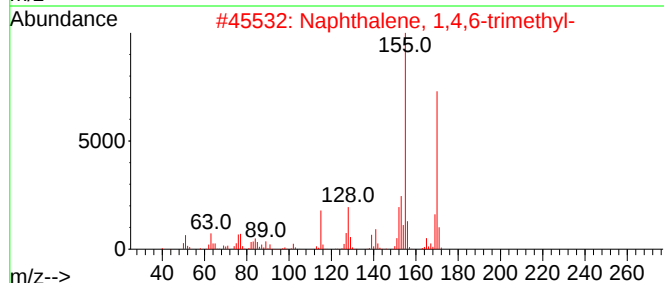
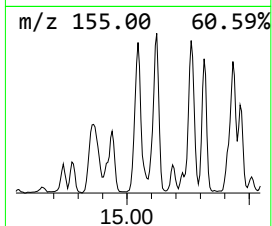
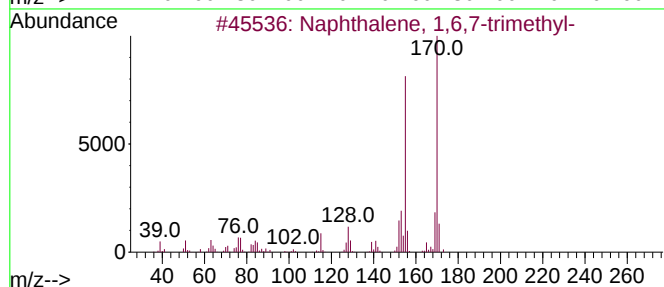
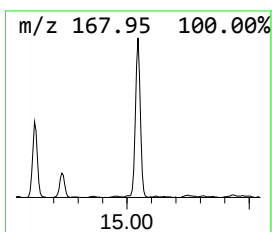
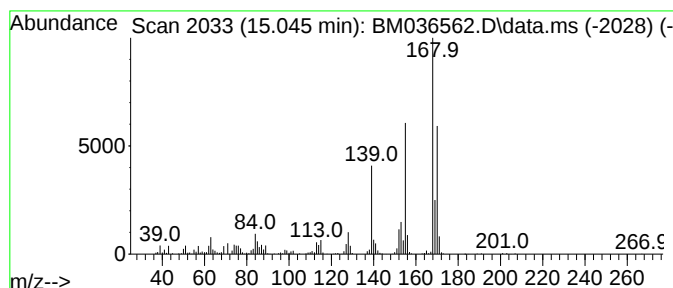
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene, 1,6,7-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.045	2.73 ng/ul	636861	Acenaphthene-d10	14.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	86
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	64
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	64
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	60
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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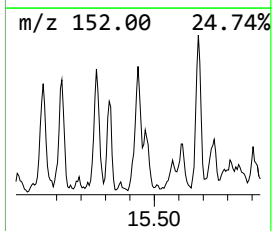
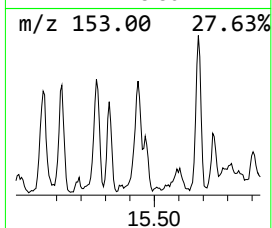
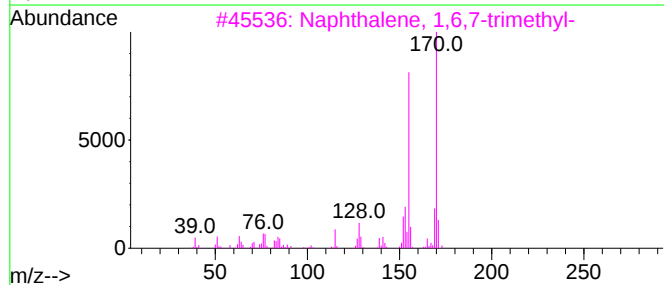
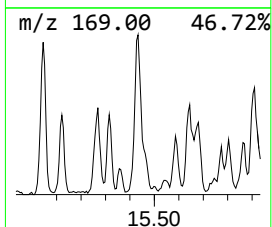
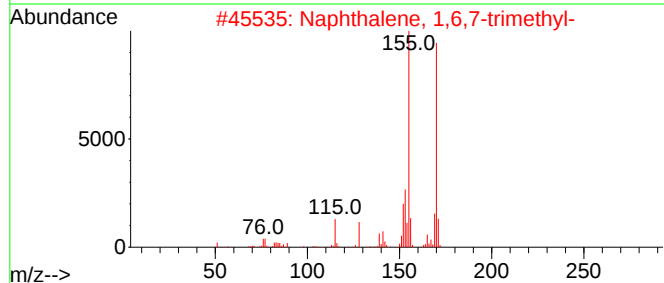
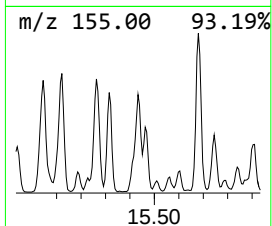
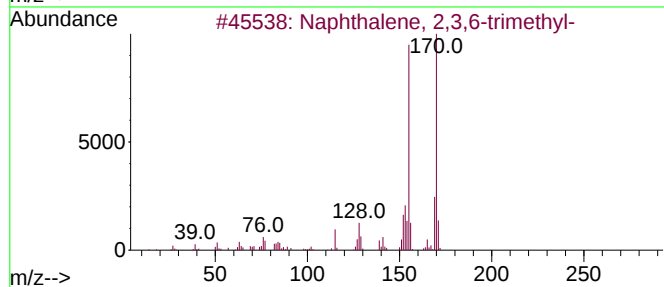
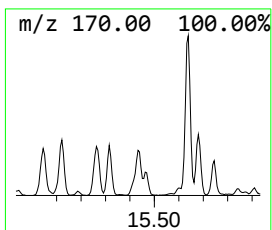
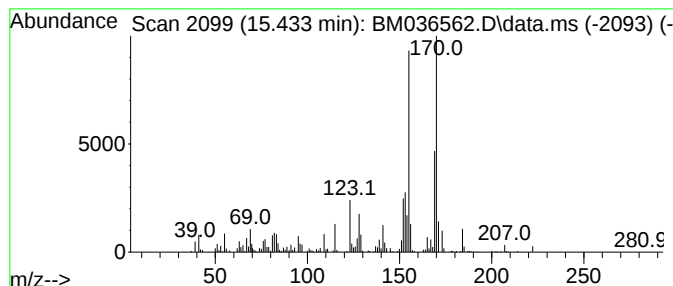
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 2,3,6-trimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	2.08 ng/ul	484803	Acenaphthene-d10	14.651

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	93
5			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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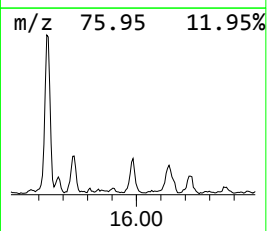
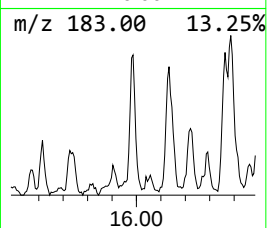
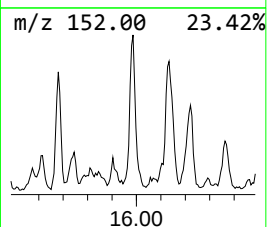
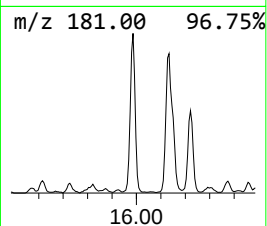
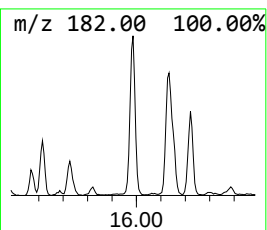
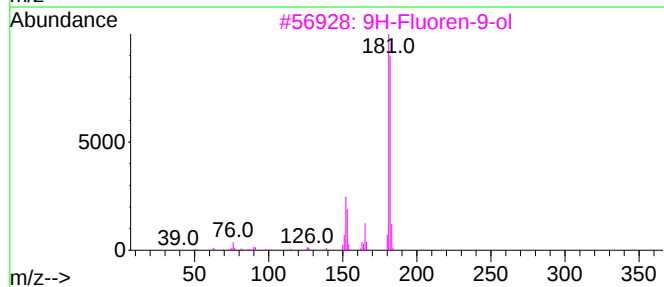
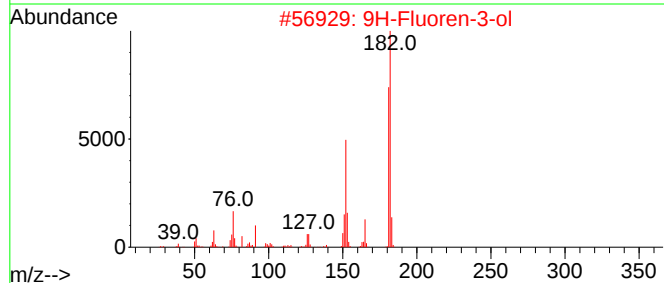
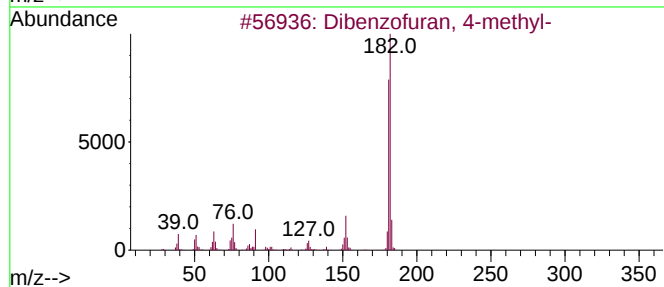
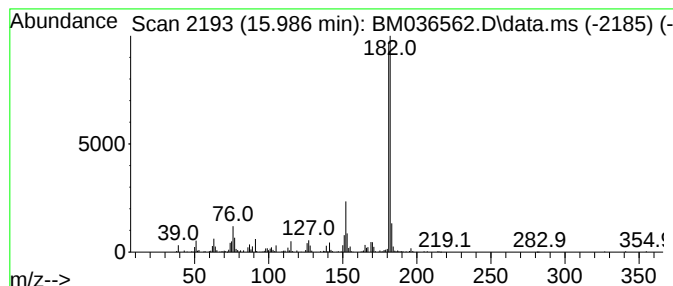
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dibenzofuran, 4-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.986	2.29 ng/ul	534419	Acenaphthene-d10	14.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2			9H-Fluoren-3-ol	182	C13H10O	006344-67-8	90
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
4			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78
5			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
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Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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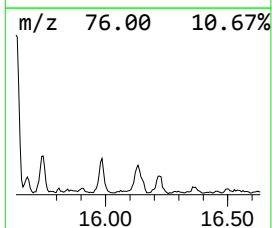
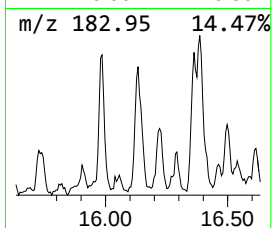
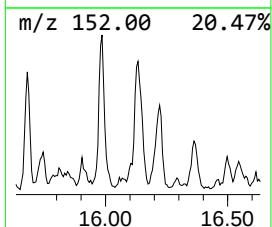
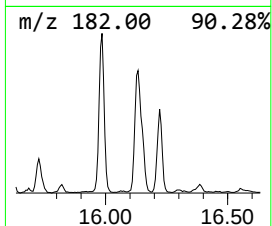
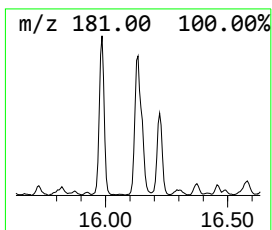
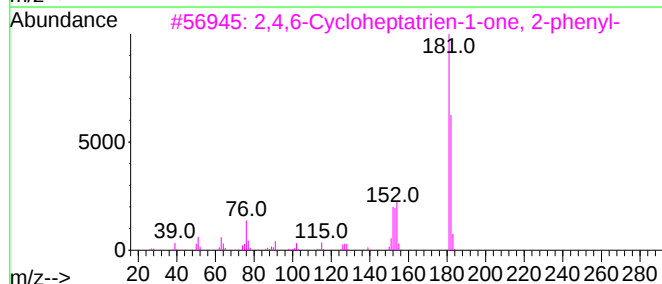
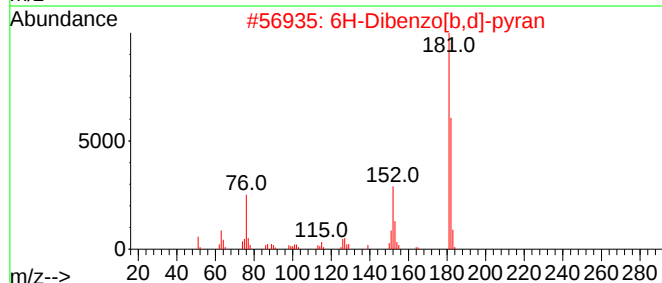
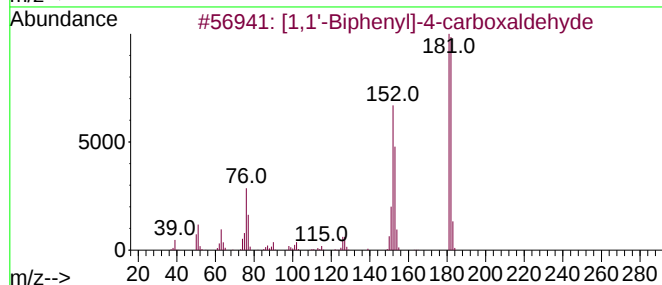
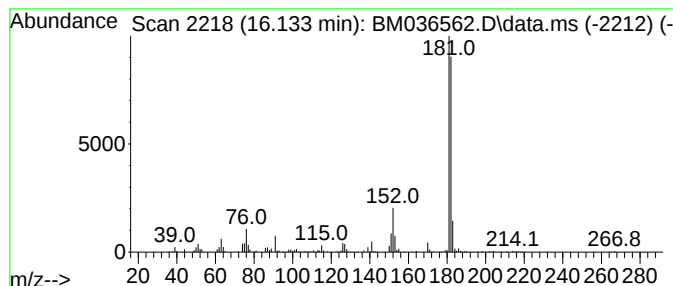
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.133	2.22 ng/ul	573091	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
2		6H-Dibenzo[b,d]-pyran	182	C13H10O	000229-95-8	87
3		2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	83
4		2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	80
5		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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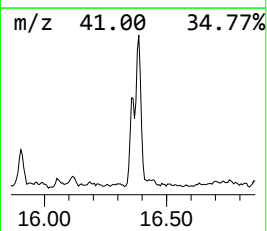
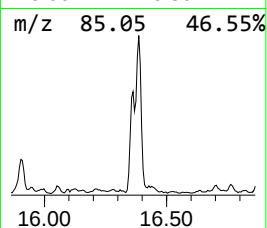
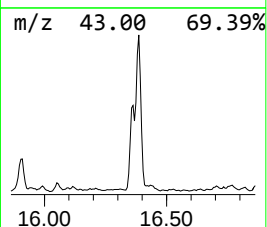
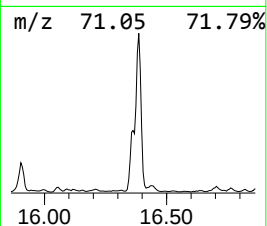
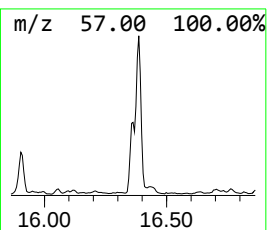
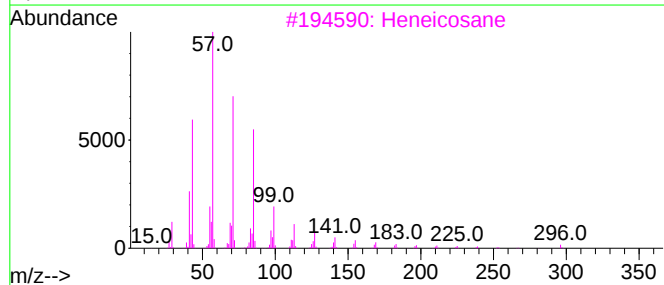
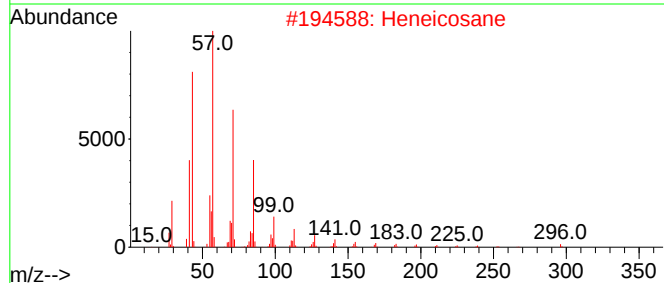
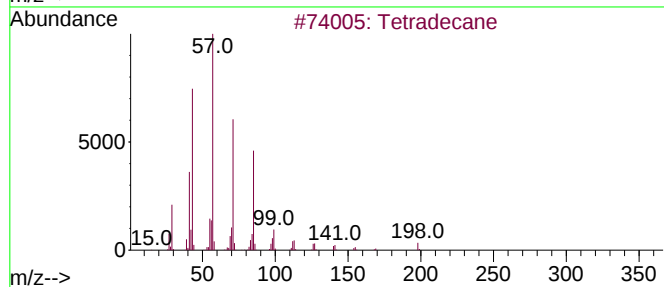
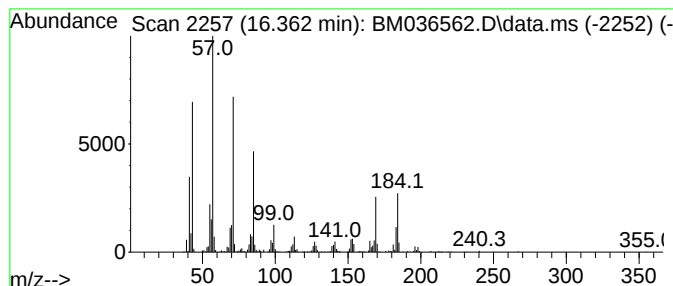
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.363	2.45 ng/ul	634050	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	64
2		Heneicosane	296	C21H44	000629-94-7	58
3		Heneicosane	296	C21H44	000629-94-7	58
4		Dodecane	170	C12H26	000112-40-3	58
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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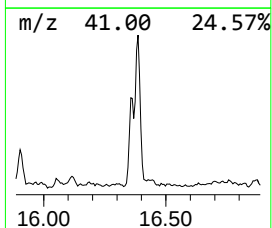
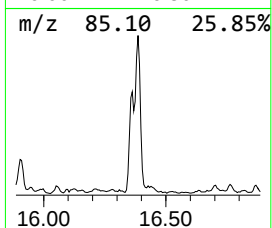
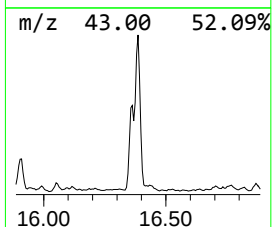
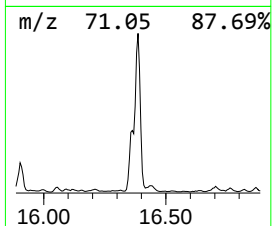
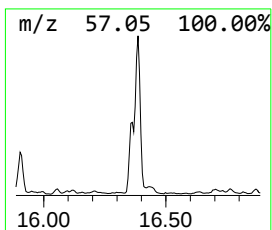
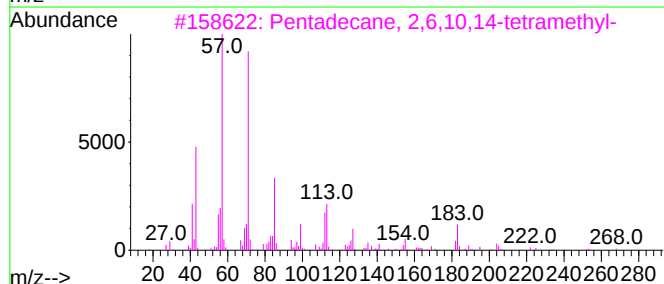
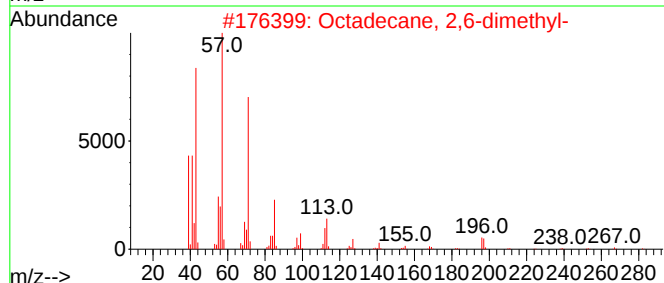
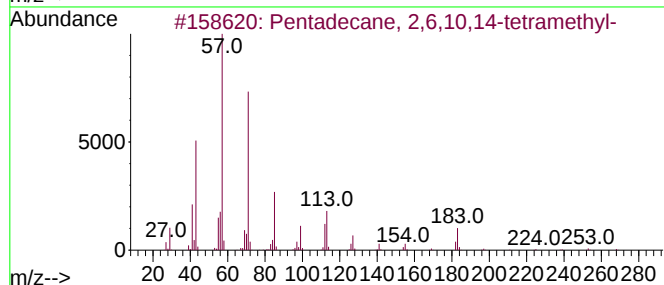
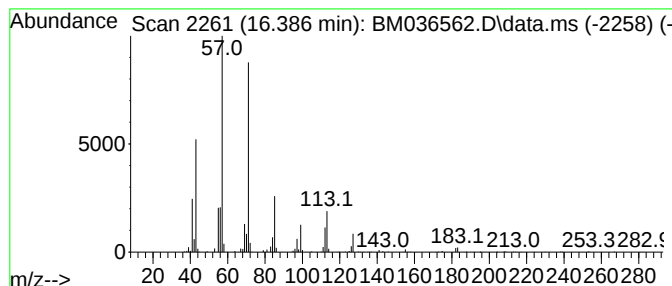
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.386	4.50 ng/ul	1163130	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2		Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	90
3		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	83
5		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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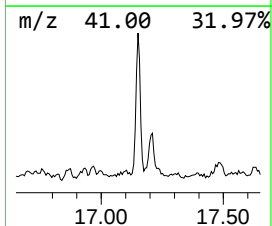
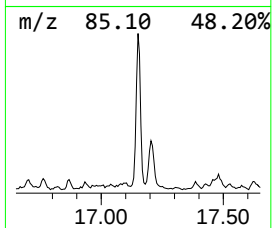
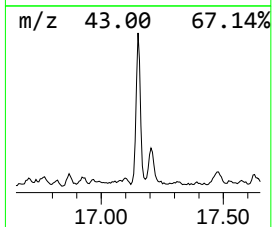
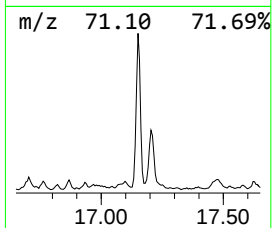
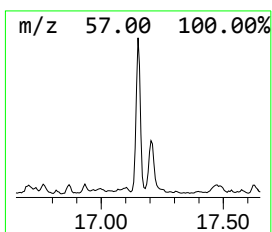
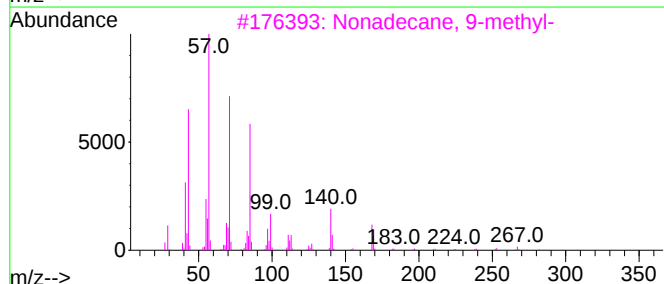
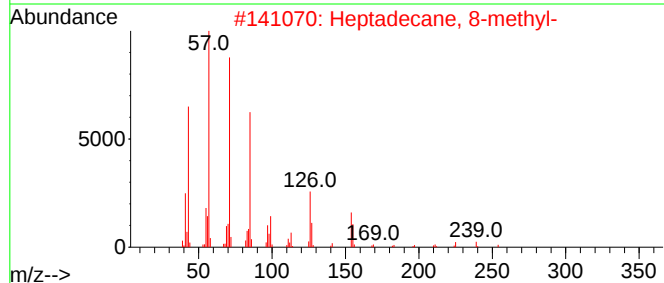
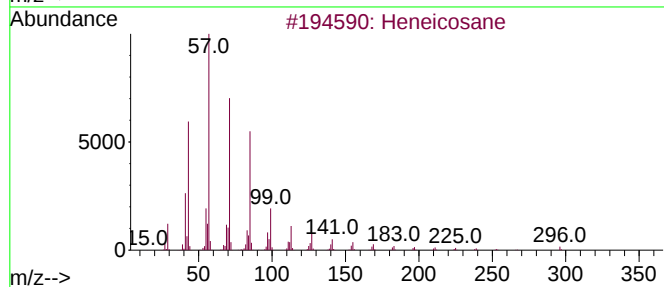
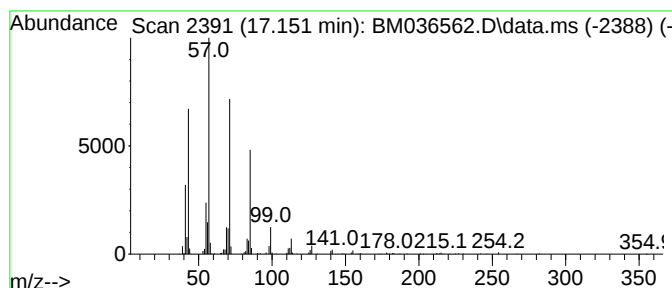
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.151	2.54 ng/ul	655837	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane, 8-methyl-	254	C18H38	013287-23-5	90
3		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		9-methylheptadecane	254	C18H38	026741-18-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
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Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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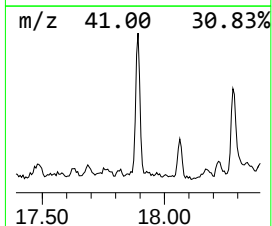
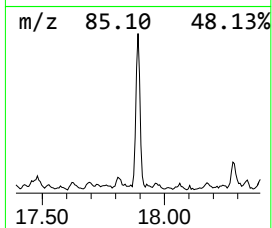
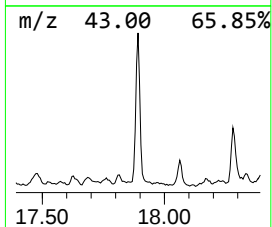
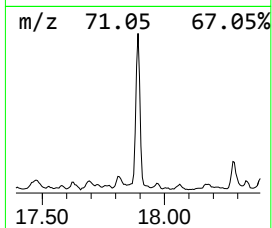
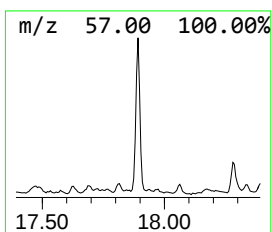
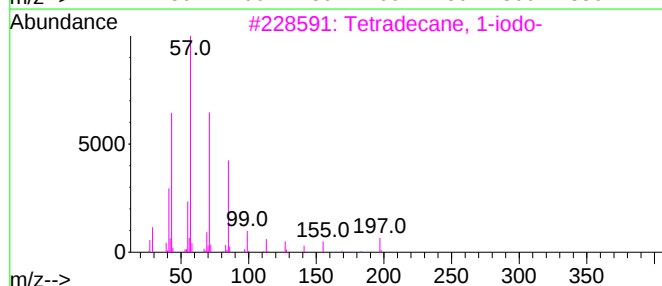
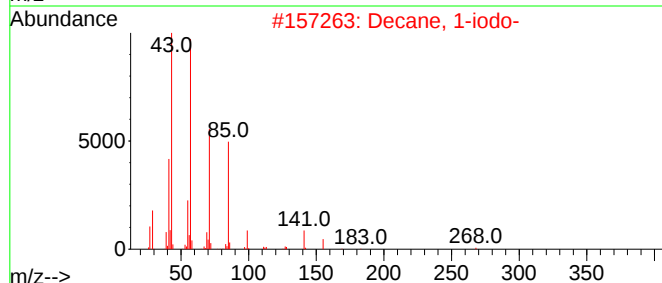
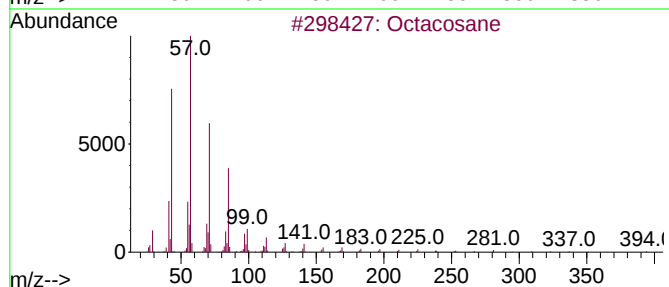
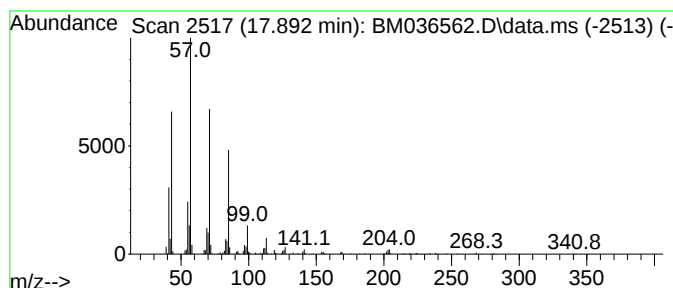
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.892	2.19 ng/ul	564930	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Decane, 1-iodo-	268	C10H21I	002050-77-3	87
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	80
4			Pentadecane	212	C15H32	000629-62-9	72
5			Nonadecane	268	C19H40	000629-92-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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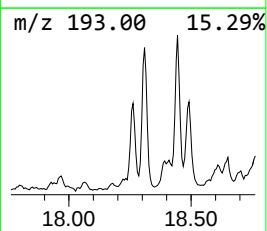
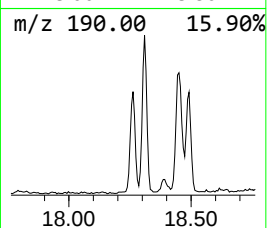
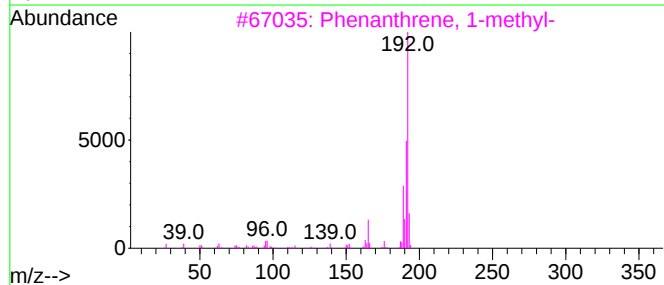
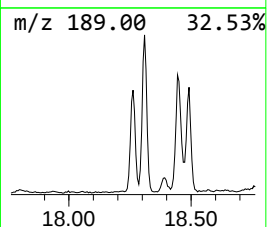
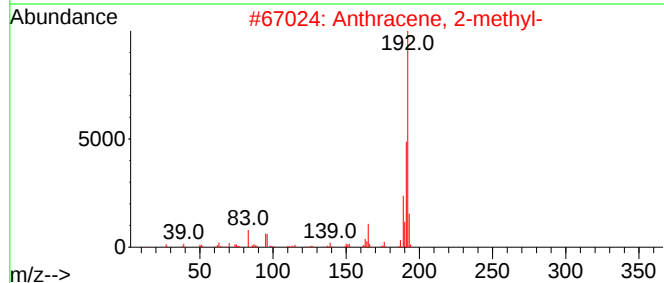
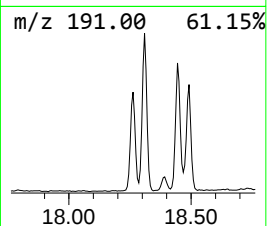
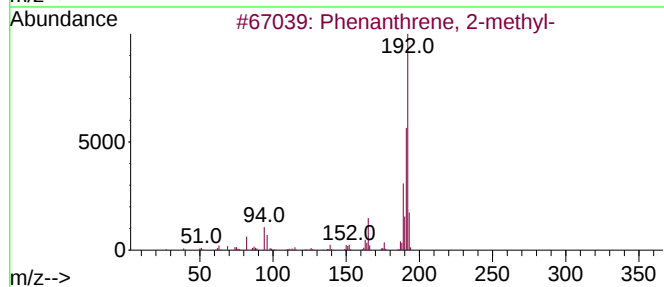
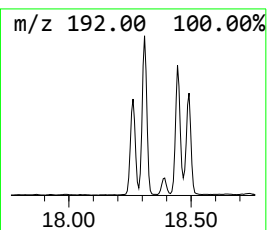
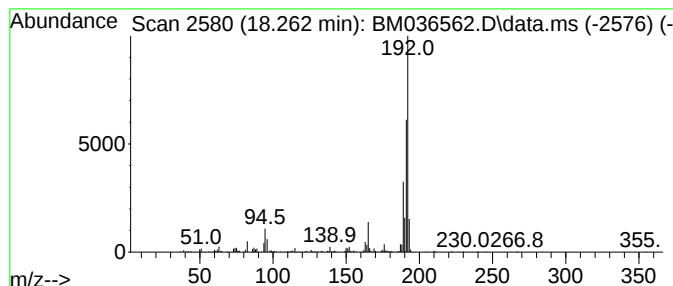
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Phenanthrene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.262	2.48 ng/ul	640574	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
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Misc :
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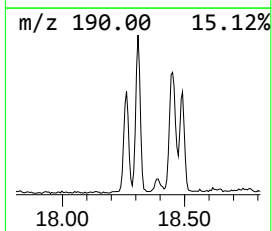
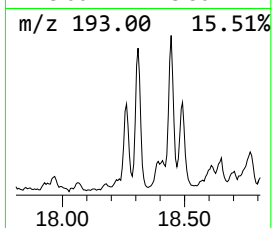
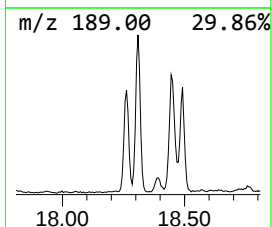
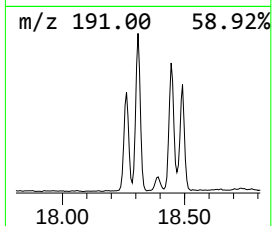
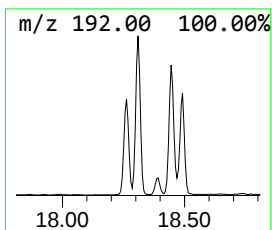
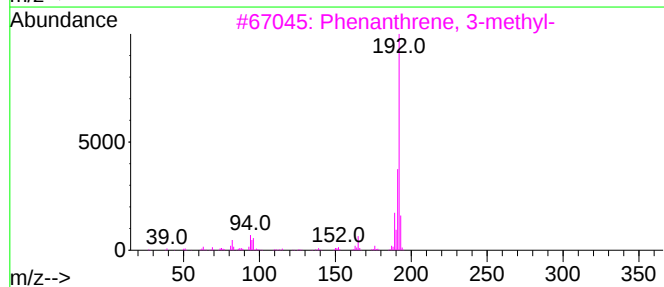
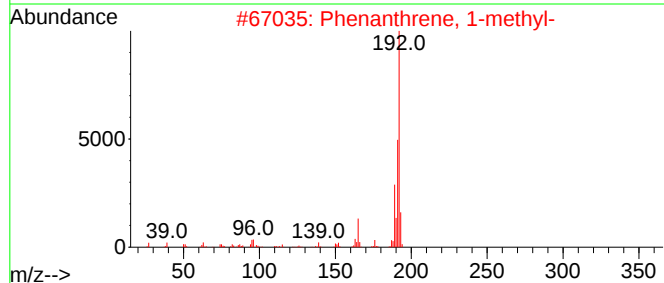
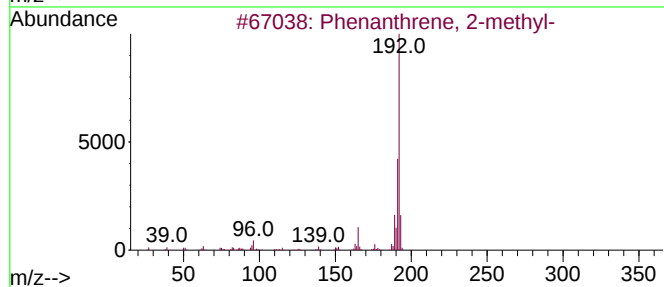
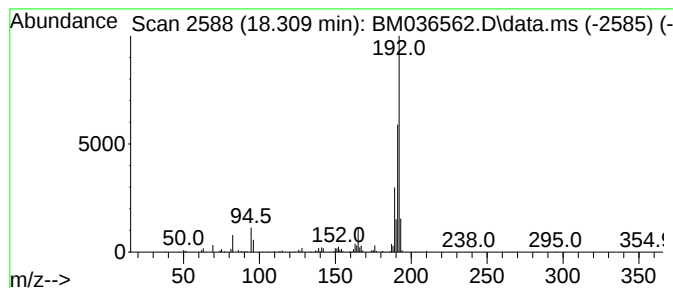
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.309	4.34 ng/ul	1122320	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	93
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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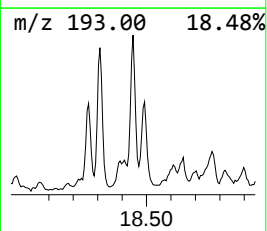
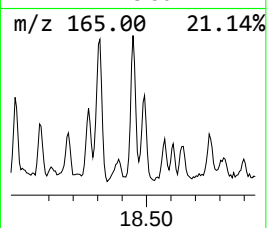
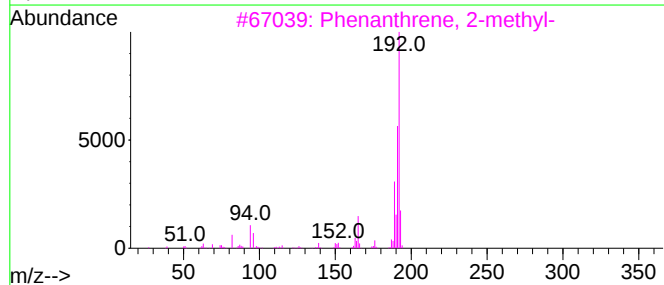
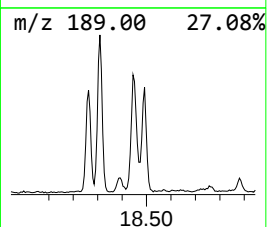
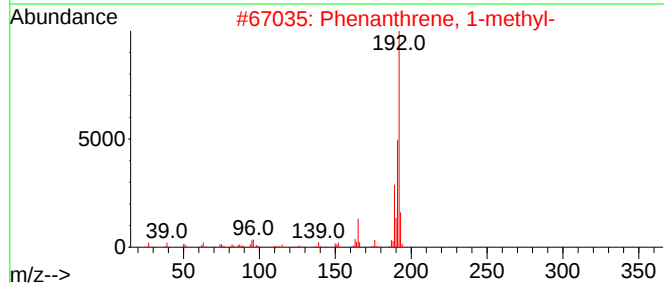
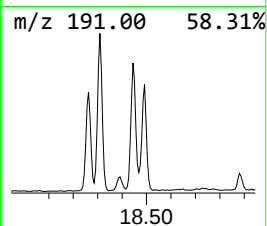
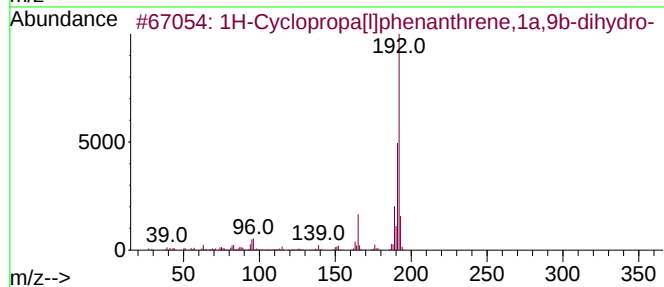
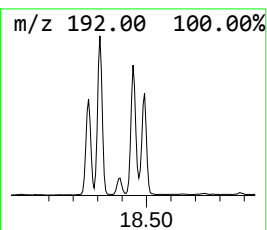
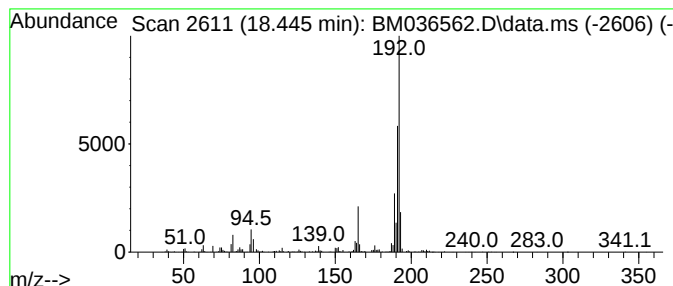
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1H-Cyclopropa[l]phenanthren... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.445	3.73 ng/ul	963228	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

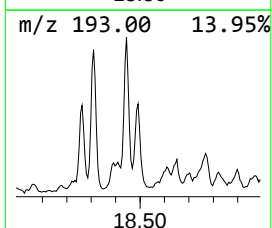
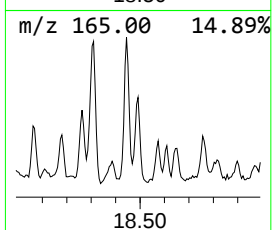
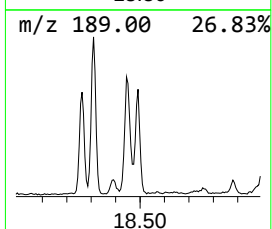
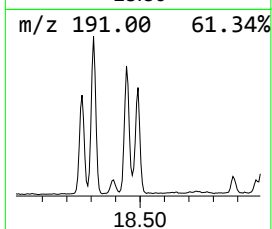
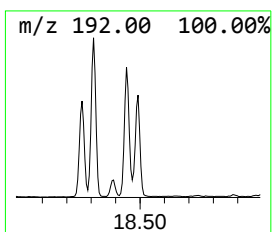
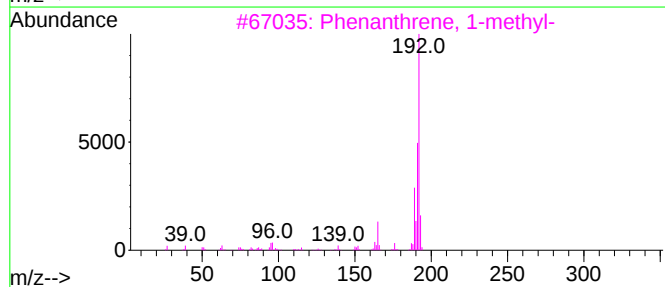
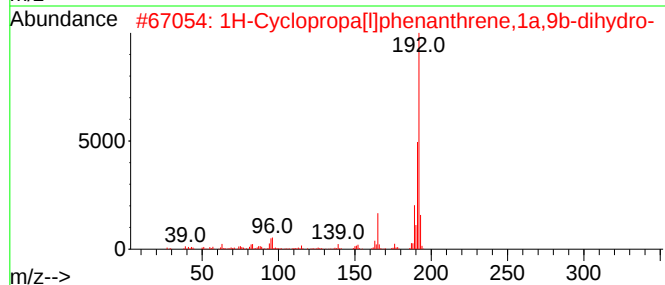
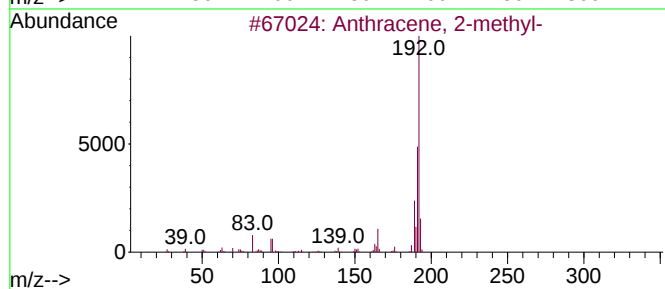
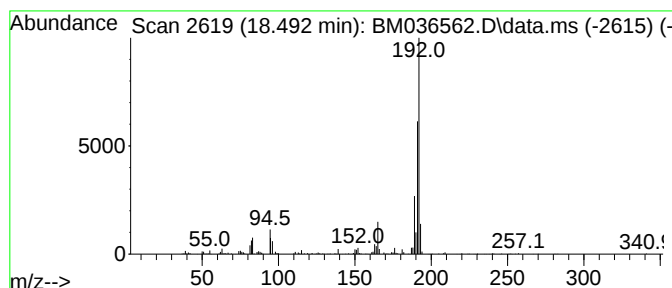
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Anthracene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.492	2.41 ng/ul	622405	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

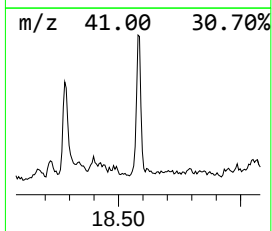
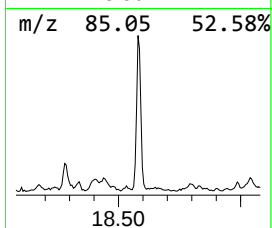
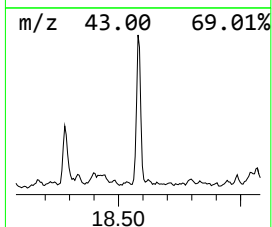
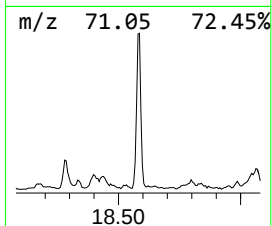
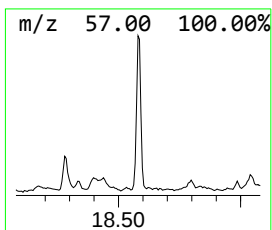
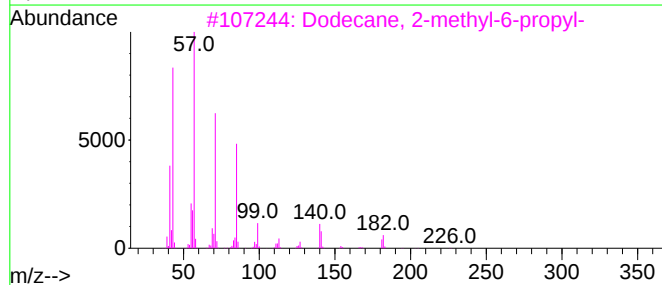
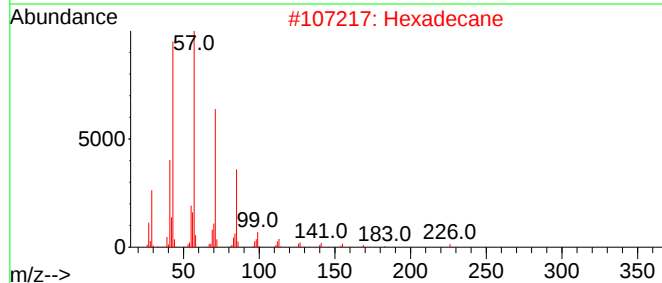
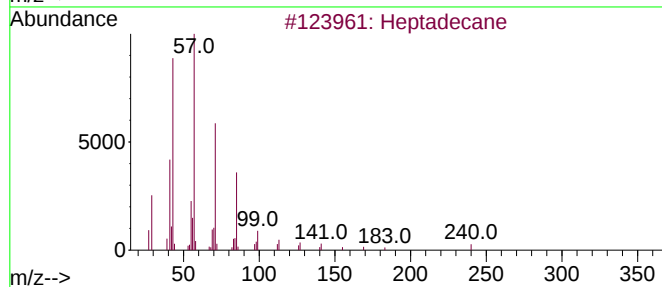
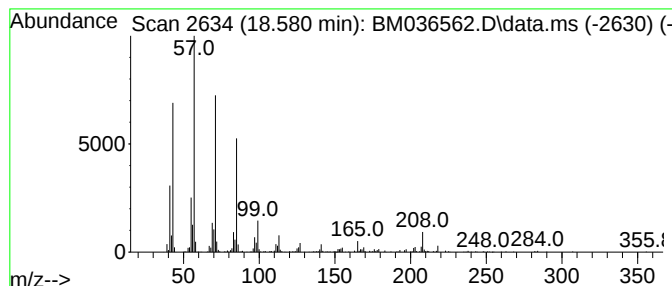
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.580	2.56 ng/ul	660036	Phenanthrene-d10		17.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	94
2	Hexadecane		226	C16H34	000544-76-3	93
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	87
4	Heneicosane		296	C21H44	000629-94-7	83
5	Nonadecane		268	C19H40	000629-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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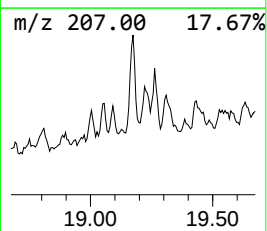
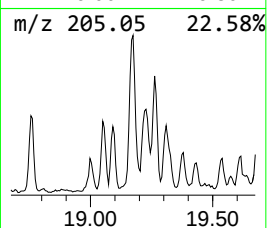
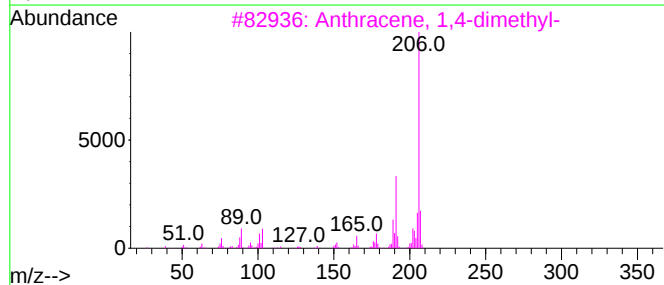
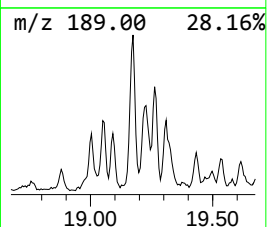
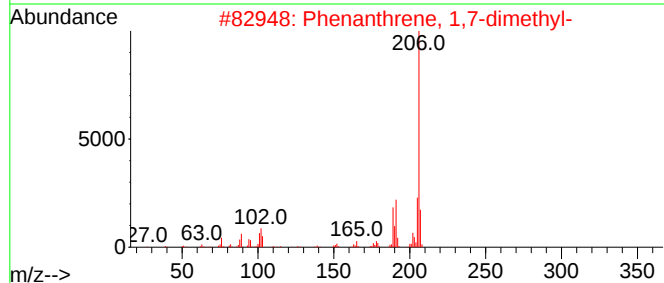
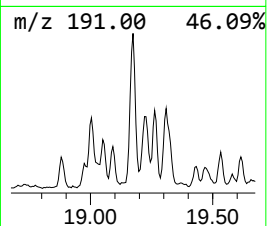
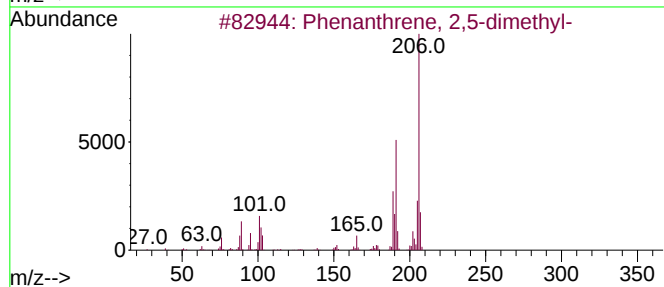
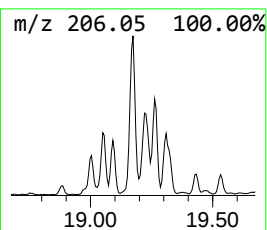
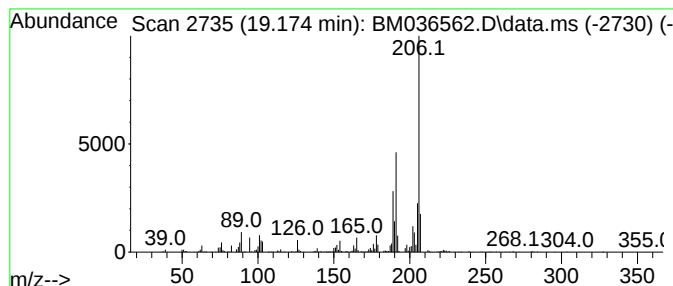
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.174	3.16 ng/ul	815574	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
3		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	94
4		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94
5		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
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Sample : N4601-14
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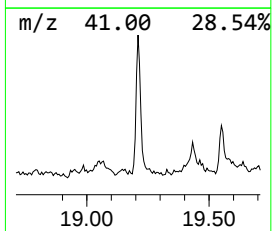
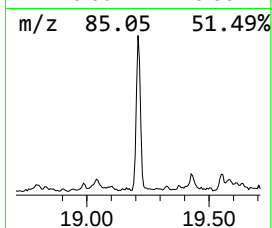
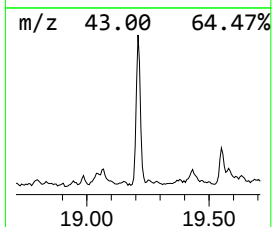
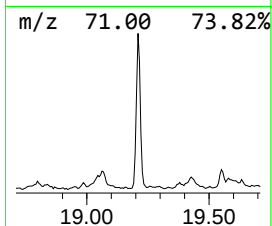
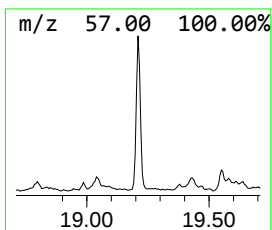
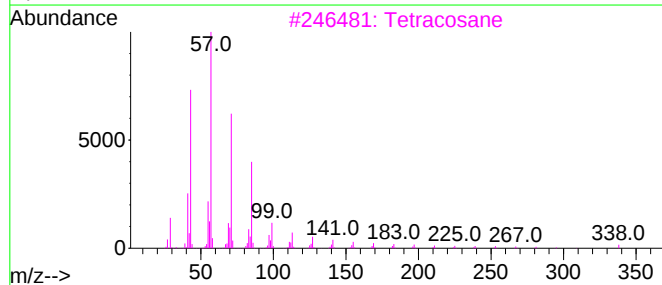
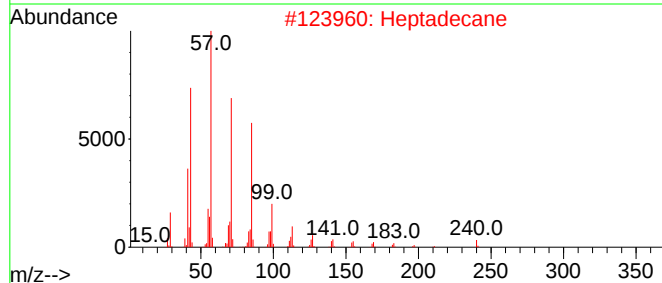
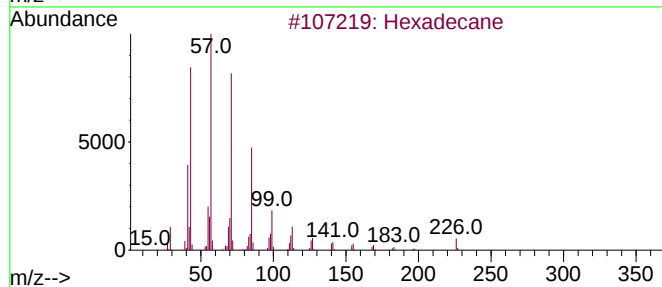
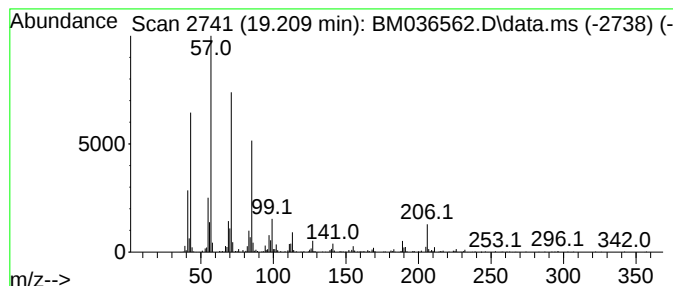
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.209	3.78 ng/ul	975654	Phenanthrene-d10	17.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Tetracosane	338	C24H50	000646-31-1	87
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
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Sample : N4601-14
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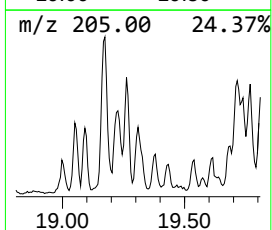
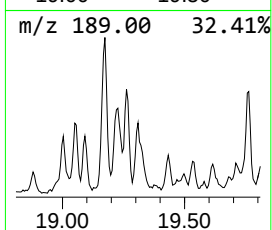
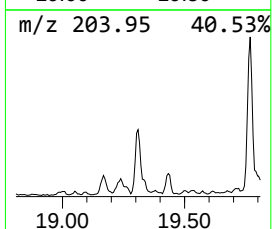
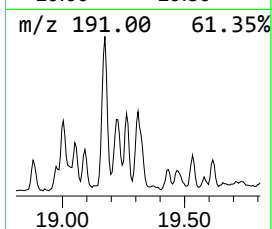
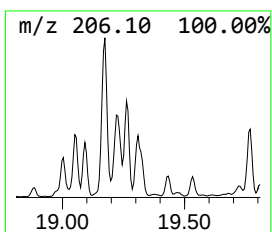
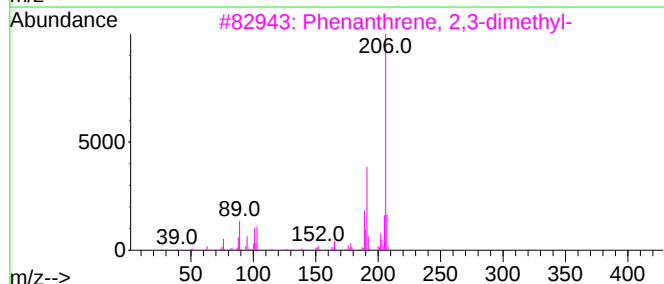
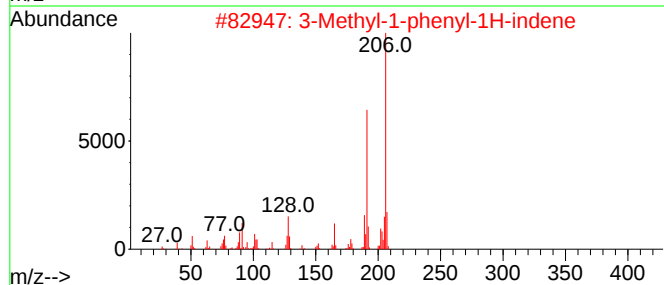
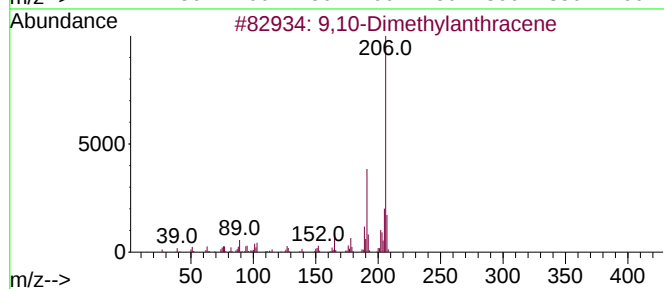
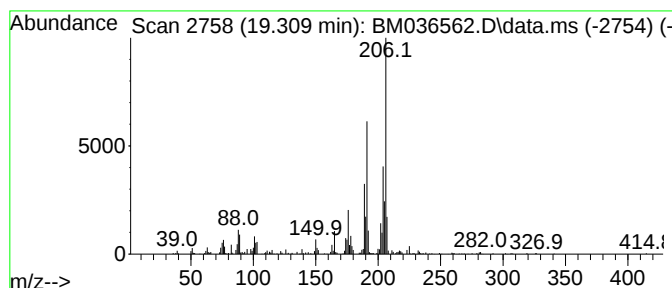
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 9,10-Dimethylantracene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.309	2.21 ng/ul	571925	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2	3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	91
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	89
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	86
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
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Sample : N4601-14
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ClientSampleId :
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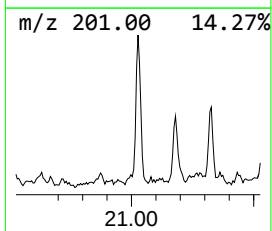
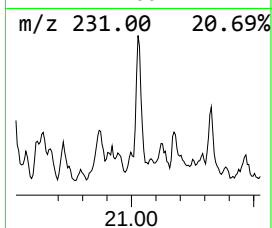
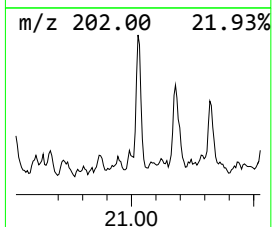
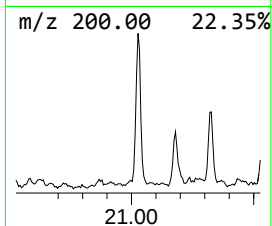
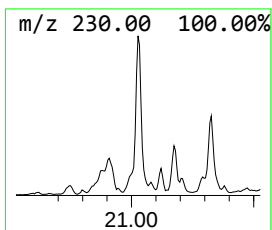
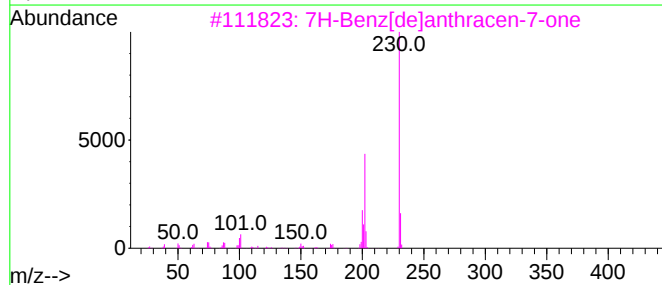
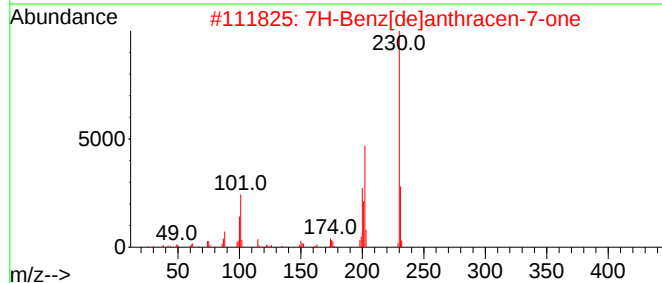
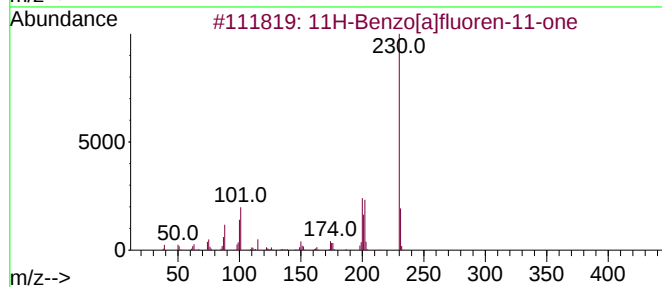
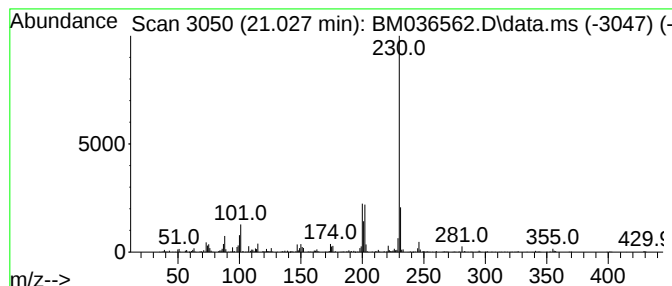
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 11H-Benzo[a]fluoren-11-one Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	2.12 ng/ul	710968	Chrysene-d12	21.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	76
4			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	72
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A5755

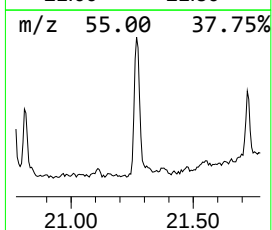
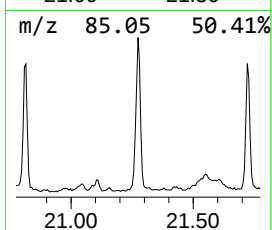
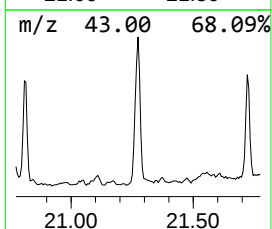
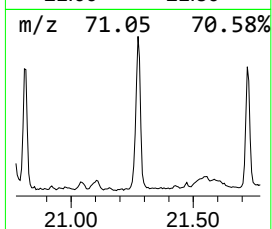
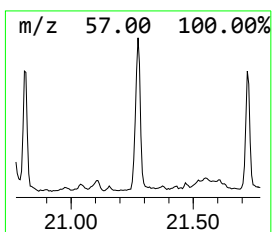
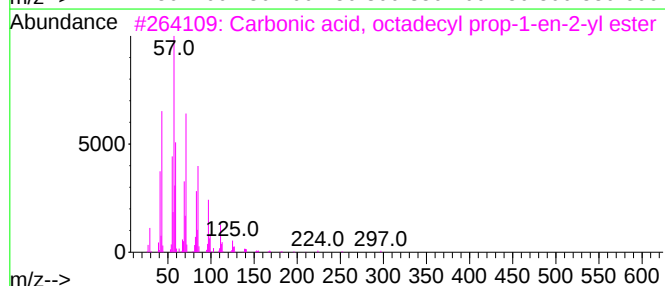
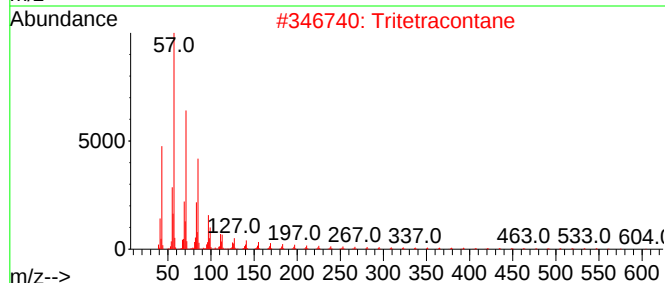
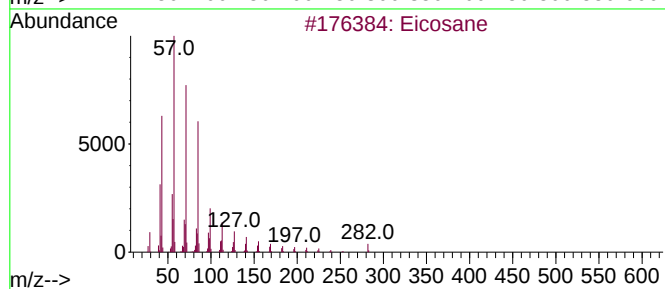
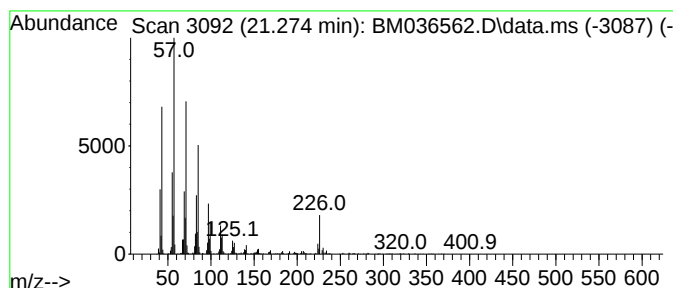
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.274	3.32 ng/ul	1115170	Chrysene-d12	21.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Tritetracontane	605	C43H88	007098-21-7	90
3			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	90
5			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

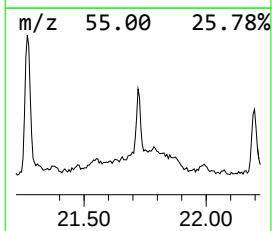
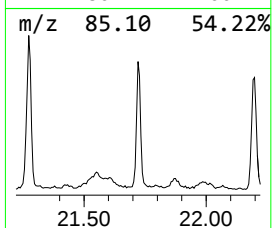
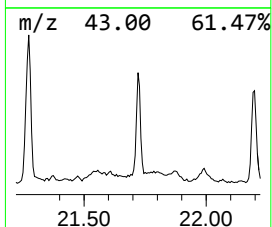
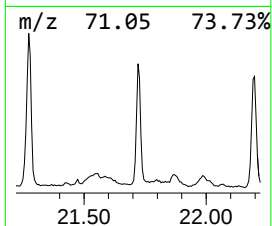
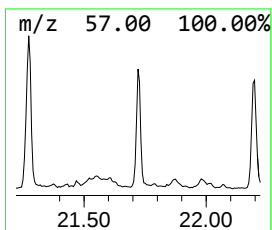
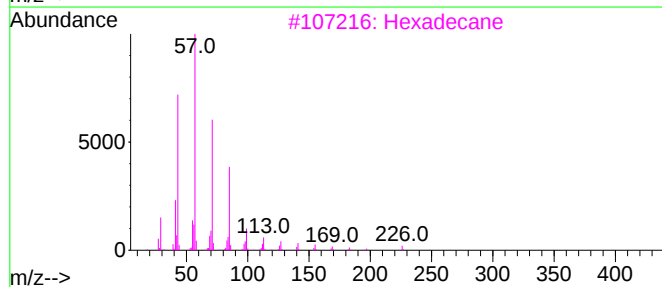
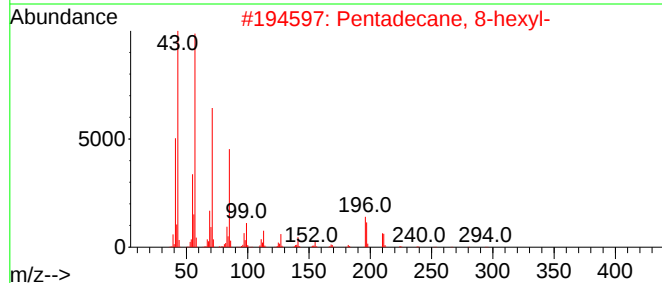
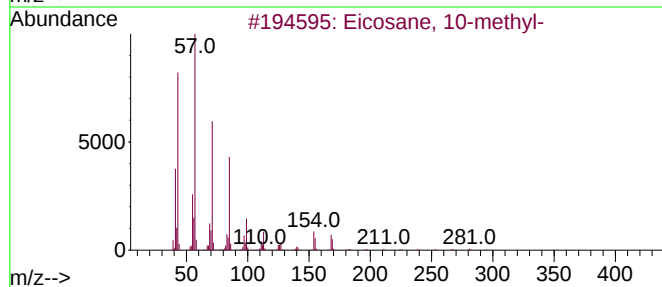
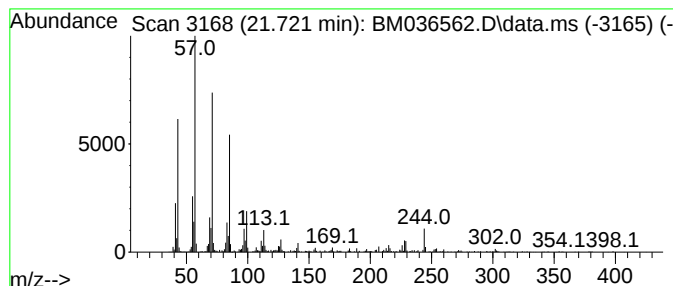
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.721	2.05 ng/ul	687297	Chrysene-d12		21.550	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-		296	C21H44	054833-23-7	93
2	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	93
3	Hexadecane		226	C16H34	000544-76-3	91
4	Octadecane		254	C18H38	000593-45-3	90
5	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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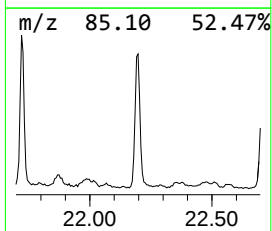
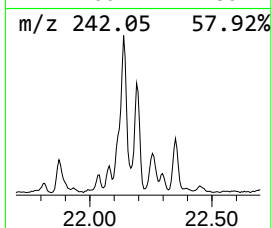
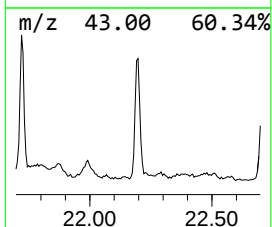
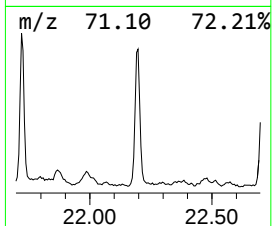
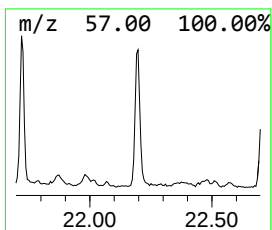
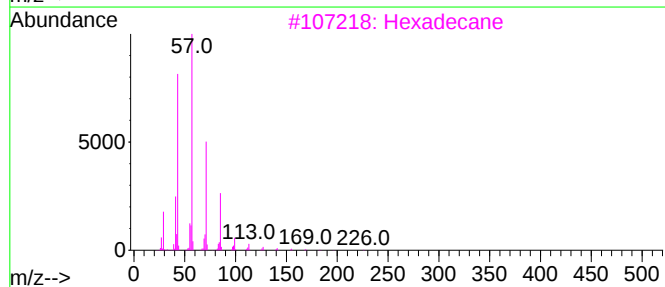
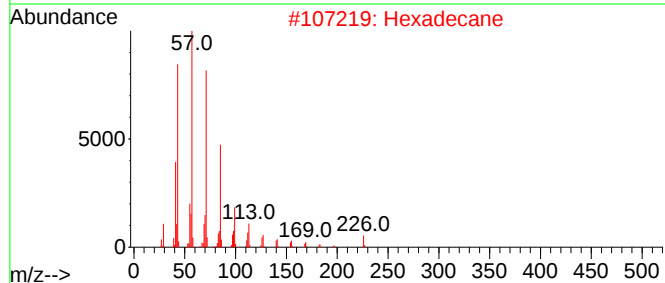
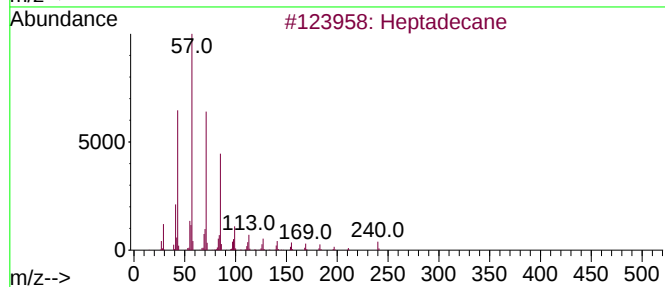
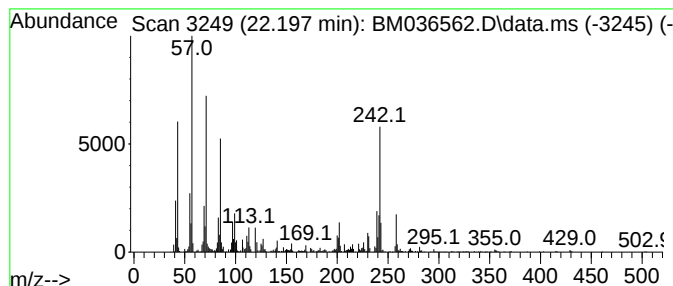
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.197	2.89 ng/ul	970472	Chrysene-d12	21.550

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Hexadecane	226	C16H34	000544-76-3	92
3		Hexadecane	226	C16H34	000544-76-3	89
4		Hexadecane	226	C16H34	000544-76-3	89
5		Heptadecane	240	C17H36	000629-78-7	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036562.D
Acq On : 16 Sep 2022 14:46
Operator : CG/JU
Sample : N4601-14
Misc :
ALS Vial : 8 Sample Multiplier: 1

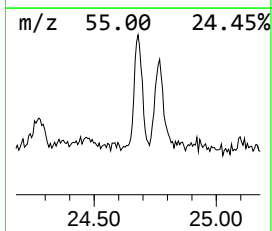
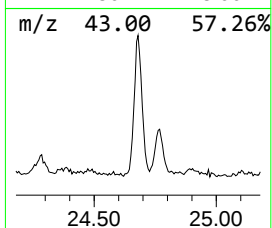
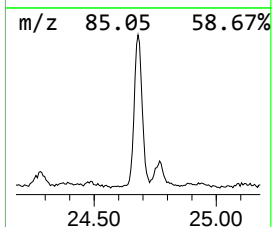
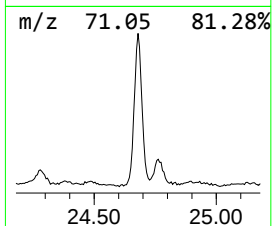
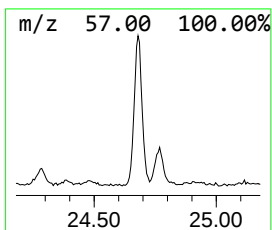
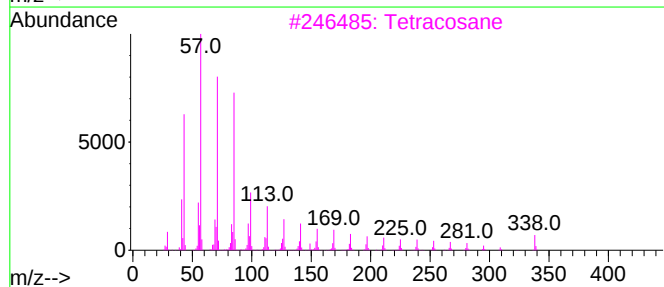
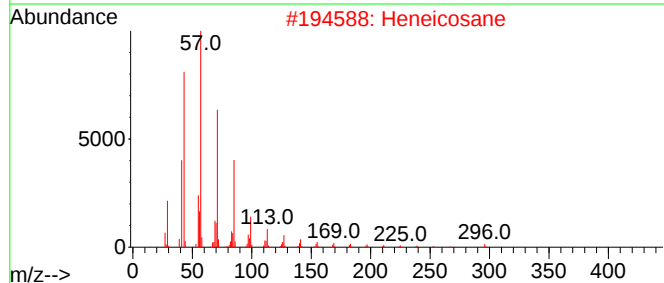
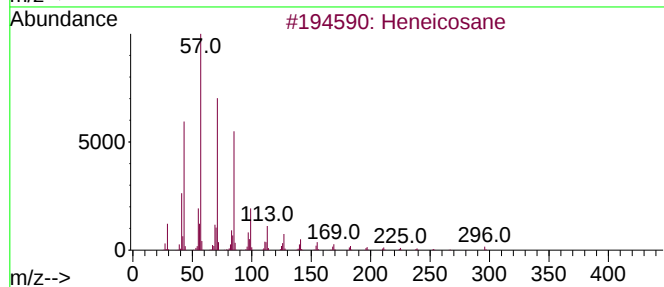
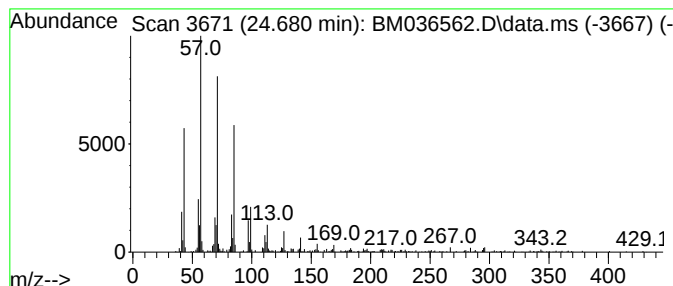
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.680	2.07 ng/ul	553376	Perylene-d12	23.997	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	98
2	Heneicosane	296	C21H44	000629-94-7	94
3	Tetracosane	338	C24H50	000646-31-1	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036562.D
 Acq On : 16 Sep 2022 14:46
 Operator : CG/JU
 Sample : N4601-14
 Misc :
 ALS Vial : 8 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
Butane, 2-metho...	3.122	157.7	ng/ul	16466200	1	8.028	2088110	20.0
Ethanol, 2-(2-b...	10.610	6.1	ng/ul	1007650	2	10.839	3326940	20.0
(DEL) Alkane: S...	11.869	2.3	ng/ul	378736	2	10.839	3326940	20.0
(DEL) Alkane: S...	13.486	2.4	ng/ul	546582	3	14.651	4661560	20.0
Naphthalene, 2,...	13.974	3.3	ng/ul	763638	3	14.651	4661560	20.0
Naphthalene, 1,...	14.210	2.1	ng/ul	486013	3	14.651	4661560	20.0
Naphthalene, 1,...	15.045	2.7	ng/ul	636861	3	14.651	4661560	20.0
Naphthalene, 2,...	15.433	2.1	ng/ul	484803	3	14.651	4661560	20.0
Dibenzofuran, 4...	15.986	2.3	ng/ul	534419	3	14.651	4661560	20.0
[1,1'-Biphenyl]...	16.133	2.2	ng/ul	573091	4	17.386	5166510	20.0
(DEL) Alkane: S...	16.363	2.5	ng/ul	634050	4	17.386	5166510	20.0
(DEL) Alkane: S...	16.386	4.5	ng/ul	1163130	4	17.386	5166510	20.0
(DEL) Alkane: S...	17.151	2.5	ng/ul	655837	4	17.386	5166510	20.0
(DEL) Alkane: S...	17.892	2.2	ng/ul	564930	4	17.386	5166510	20.0
Phenanthrene, 2...	18.262	2.5	ng/ul	640574	4	17.386	5166510	20.0
Phenanthrene, 1...	18.309	4.3	ng/ul	1122320	4	17.386	5166510	20.0
1H-Cyclopropa[l...	18.445	3.7	ng/ul	963228	4	17.386	5166510	20.0
Anthracene, 2-m...	18.492	2.4	ng/ul	622405	4	17.386	5166510	20.0
(DEL) Alkane: S...	18.580	2.6	ng/ul	660036	4	17.386	5166510	20.0
Phenanthrene, 2...	19.174	3.2	ng/ul	815574	4	17.386	5166510	20.0
(DEL) Alkane: S...	19.209	3.8	ng/ul	975654	4	17.386	5166510	20.0
9,10-Dimethylan...	19.309	2.2	ng/ul	571925	4	17.386	5166510	20.0
11H-Benzo[a]flu...	21.027	2.1	ng/ul	710968	5	21.550	6721310	20.0
(DEL) Alkane: S...	21.274	3.3	ng/ul	1115170	5	21.550	6721310	20.0
(DEL) Alkane: S...	21.721	2.0	ng/ul	687297	5	21.550	6721310	20.0
(DEL) Alkane: S...	22.197	2.9	ng/ul	970472	5	21.550	6721310	20.0
(DEL) Alkane: S...	24.680	2.1	ng/ul	553376	6	23.997	5344650	20.0