

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
 Data File : BP020772.D
 Acq On : 03 Jul 2024 15:56
 Operator : MA/JU
 Sample : P2963-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CR7

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_P\Methods\SFAM-EPA-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020772.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.546	91	95	104	rBV	51735	80987	1.06%	0.068%
2	3.999	167	172	181	rVB	70543	105461	1.38%	0.088%
3	4.899	319	325	333	rBV	140905	212468	2.78%	0.178%
4	6.928	661	670	687	rVB	958850	1609068	21.07%	1.349%
5	7.087	690	697	709	rVB	771686	1158208	15.17%	0.971%
6	7.281	722	730	738	rBV	952525	1598595	20.94%	1.340%
7	7.740	799	808	815	rBV	699478	1232739	16.14%	1.033%
8	8.446	918	928	936	rBV	1145631	1999066	26.18%	1.676%
9	8.552	942	946	955	rVB	87289	150797	1.97%	0.126%
10	8.887	995	1003	1015	rBV	995476	1579347	20.68%	1.324%
11	9.434	1088	1096	1104	rBV2	95468	159934	2.09%	0.134%
12	9.599	1115	1124	1135	rBV	839851	1387051	18.17%	1.163%
13	10.134	1207	1215	1235	rBV	1203673	2114782	27.70%	1.773%
14	10.499	1269	1277	1282	rBV	1034396	1775695	23.26%	1.489%
15	10.546	1282	1285	1292	rVB	125058	207209	2.71%	0.174%
16	10.646	1295	1302	1321	rBV	252499	562010	7.36%	0.471%
17	11.034	1361	1368	1377	rBV3	43770	85422	1.12%	0.072%
18	11.240	1396	1403	1409	rBV	118385	229062	3.00%	0.192%
19	12.157	1553	1559	1567	rVB	155448	231289	3.03%	0.194%
20	12.387	1591	1598	1605	rVB	130833	221131	2.90%	0.185%
21	13.181	1726	1733	1740	rVB2	44660	85587	1.12%	0.072%
22	13.381	1761	1767	1776	rVB	55630	130481	1.71%	0.109%
23	13.534	1778	1793	1798	rBV3	81274	214315	2.81%	0.180%
24	13.669	1809	1816	1821	rBV	149321	273990	3.59%	0.230%
25	13.769	1821	1833	1847	rVB	1829742	3175154	41.58%	2.662%
26	13.910	1852	1857	1861	rBV	90644	127047	1.66%	0.106%
27	14.045	1869	1880	1897	rBV	2362451	3978821	52.11%	3.335%
28	14.351	1922	1932	1939	rBV	1571739	2507607	32.84%	2.102%
29	14.422	1939	1944	1951	rVB	377195	607950	7.96%	0.510%
30	14.593	1963	1973	1982	rBV2	889202	1778374	23.29%	1.491%
31	14.675	1982	1987	1991	rVV2	100642	200965	2.63%	0.168%
32	14.763	1991	2002	2010	rVV2	287154	653637	8.56%	0.548%
33	14.834	2010	2014	2018	rVB	78543	114731	1.50%	0.096%
34	14.987	2033	2040	2045	rBV2	71129	141894	1.86%	0.119%
35	15.040	2045	2049	2054	rVB2	73376	104293	1.37%	0.087%
36	15.157	2062	2069	2072	rBV3	76315	150942	1.98%	0.127%

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Integrator: RTE

Smoothing : OFF

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Title : SVOA CALIBRATION

37	15.369	2098	2105	2110	rVV	3006905	4521669	59.22%	3.790%
38	15.422	2110	2114	2120	rVV	463626	725326	9.50%	0.608%
39	15.504	2123	2128	2133	rVV	806573	1154206	15.12%	0.968%
40	15.545	2133	2135	2138	rVV	150025	184299	2.41%	0.154%

41	15.592	2138	2143	2147	rVV4	110036	194665	2.55%	0.163%
42	15.640	2147	2151	2154	rVV2	84815	115656	1.51%	0.097%
43	15.681	2154	2158	2161	rVV3	51355	93331	1.22%	0.078%
44	15.728	2162	2166	2174	rVB2	133074	246989	3.23%	0.207%
45	15.869	2187	2190	2198	rVV	119298	265147	3.47%	0.222%

46	15.963	2198	2206	2213	rVB3	48323	138354	1.81%	0.116%
47	16.116	2228	2232	2237	rVB4	111735	189007	2.48%	0.158%
48	16.434	2279	2286	2287	rBV2	166130	261413	3.42%	0.219%
49	16.451	2287	2289	2293	rVV2	195350	254194	3.33%	0.213%
50	16.498	2293	2297	2303	rVB2	181306	264622	3.47%	0.222%

51	16.551	2303	2306	2311	rBV4	61773	125220	1.64%	0.105%
52	16.598	2312	2314	2320	rVB2	113598	159267	2.09%	0.134%
53	16.669	2321	2326	2327	rBV2	68673	115526	1.51%	0.097%
54	16.728	2333	2336	2340	rVB4	90856	118207	1.55%	0.099%
55	16.816	2346	2351	2357	rVB	266137	377917	4.95%	0.317%

56	16.910	2360	2367	2373	rVV4	208841	516577	6.77%	0.433%
57	16.975	2374	2378	2388	rVB	335131	574371	7.52%	0.481%
58	17.169	2405	2411	2414	rBV	1891388	2738153	35.86%	2.295%
59	17.216	2414	2419	2424	rVV	4616474	7635607	100.00%	6.401%
60	17.275	2424	2429	2432	rVV	3326596	5059853	66.27%	4.242%

61	17.304	2432	2434	2440	rVV	1018846	1311568	17.18%	1.099%
62	17.363	2440	2444	2450	rVB3	111981	217094	2.84%	0.182%
63	17.581	2477	2481	2486	rBV2	256954	416705	5.46%	0.349%
64	17.769	2508	2513	2520	rBV2	206269	415443	5.44%	0.348%
65	17.916	2534	2538	2543	rVB	182073	295082	3.86%	0.247%

66	18.075	2559	2565	2567	rBV	848313	1390506	18.21%	1.166%
67	18.104	2567	2570	2571	rVV2	953011	1211048	15.86%	1.015%
68	18.128	2571	2574	2580	rVB	1142914	1885619	24.70%	1.581%
69	18.216	2584	2589	2593	rVB	369548	540258	7.08%	0.453%
70	18.281	2593	2600	2603	rBV2	1342791	2640257	34.58%	2.213%

71	18.310	2603	2605	2613	rVB	954422	1217267	15.94%	1.020%
72	18.598	2649	2654	2657	rBV	524121	783976	10.27%	0.657%
73	18.645	2658	2662	2666	rVV2	386238	633236	8.29%	0.531%
74	18.686	2666	2669	2672	rVV4	151209	222715	2.92%	0.187%
75	18.722	2673	2675	2681	rVB	159382	238965	3.13%	0.200%

76	18.845	2693	2696	2701	rVB	318309	402861	5.28%	0.338%
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 Title : SVOA CALIBRATION

77	18.898	2703	2705	2709	rVB	189868	221549	2.90%	0.186%
78	19.028	2723	2727	2732	rBV	613155	1028309	13.47%	0.862%
79	19.092	2732	2738	2742	rVV4	302019	681386	8.92%	0.571%
80	19.175	2748	2752	2761	rBV3	445424	1130442	14.80%	0.948%
81	19.304	2768	2774	2780	rVB	4878733	7161190	93.79%	6.003%
82	19.369	2783	2785	2789	rBV2	159937	195922	2.57%	0.164%
83	19.410	2789	2792	2794	rBV2	209876	278220	3.64%	0.233%
84	19.439	2794	2797	2804	rVB4	438598	773048	10.12%	0.648%
85	19.651	2828	2833	2836	rBV2	3945616	6388081	83.66%	5.355%
86	19.686	2836	2839	2847	rVV	4151926	5620840	73.61%	4.712%
87	19.763	2848	2852	2858	rVB4	319540	601421	7.88%	0.504%
88	20.163	2916	2920	2922	rBV4	403184	628658	8.23%	0.527%
89	20.198	2923	2926	2930	rVB	854548	1084361	14.20%	0.909%
90	20.304	2941	2944	2947	rBV	514939	544222	7.13%	0.456%
91	20.357	2951	2953	2957	rVB	602253	679992	8.91%	0.570%
92	20.510	2976	2979	2983	rBV	378308	533761	6.99%	0.447%
93	20.557	2984	2987	2992	rVV	482258	746535	9.78%	0.626%
94	21.227	3098	3101	3105	rVB	456690	594909	7.79%	0.499%
95	21.280	3105	3110	3116	rBV4	1217419	2207705	28.91%	1.851%
96	21.604	3159	3165	3172	rBV3	2379000	6567650	86.01%	5.505%
97	21.669	3172	3176	3186	rVB	2018669	3469626	45.44%	2.908%
98	23.904	3550	3556	3557	rBV	1085217	1808179	23.68%	1.516%
99	24.733	3692	3697	3705	rVB2	1040470	2613855	34.23%	2.191%
100	24.957	3730	3735	3745	rVB	1133263	2935016	38.44%	2.460%

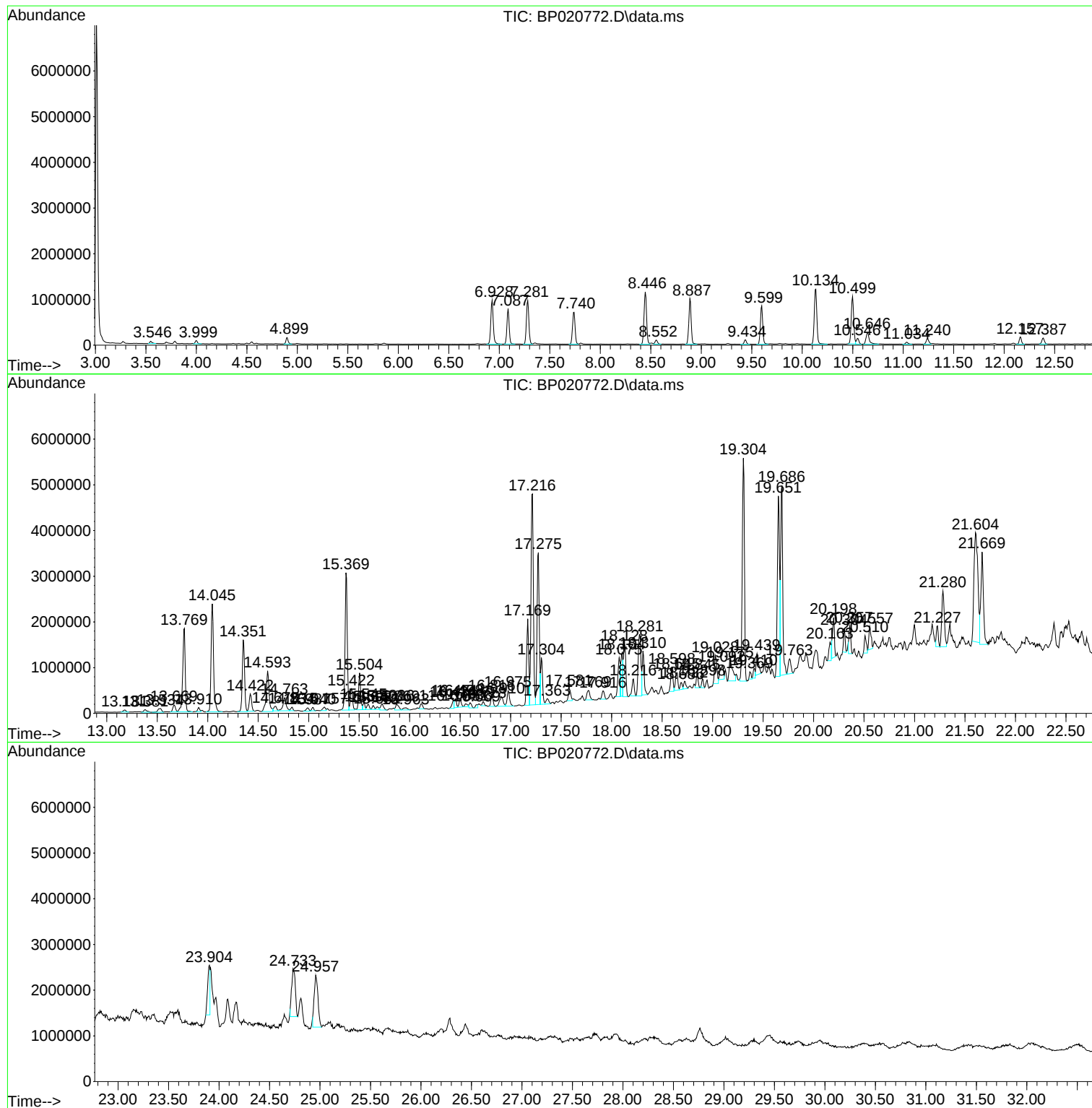
Sum of corrected areas: 119293162

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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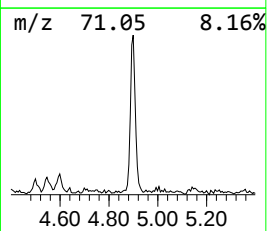
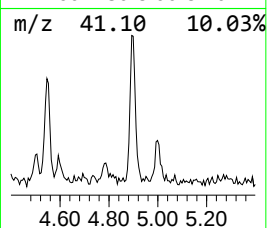
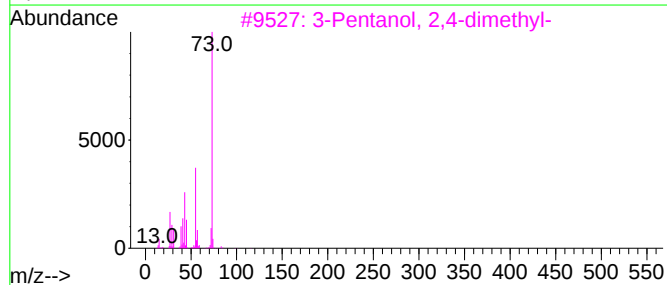
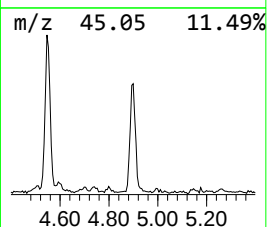
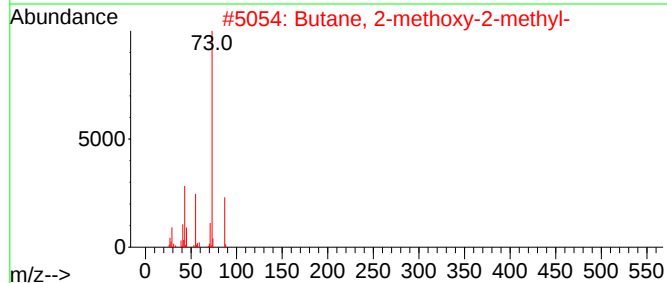
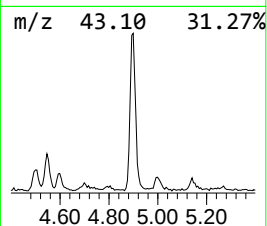
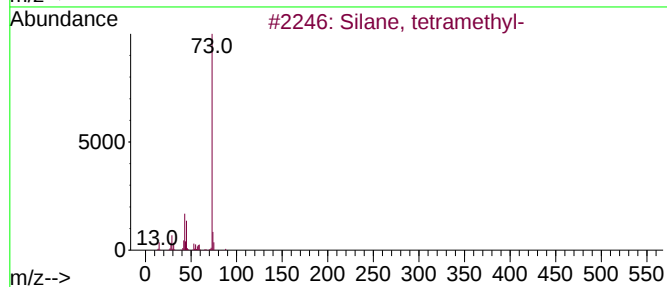
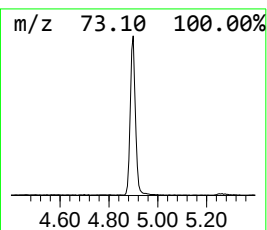
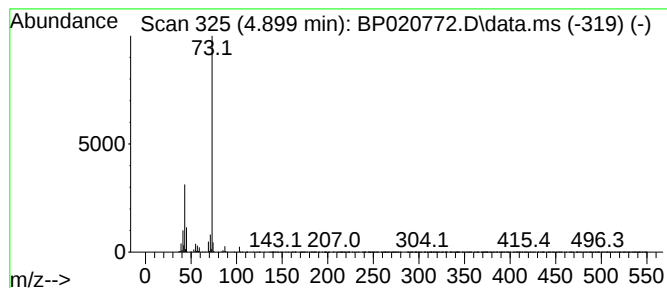
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Silane, tetramethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.899	3.45 ng/ul	212468	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, tetramethyl-	88	C4H12Si	000075-76-3	50
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	42
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	33
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
5			1,3-Dioxolane, 2-(3-bromo-3-bute...	206	C7H11BrO2	333961-98-1	9



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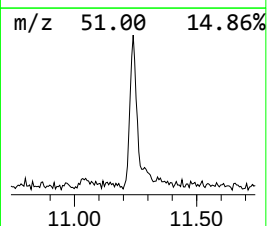
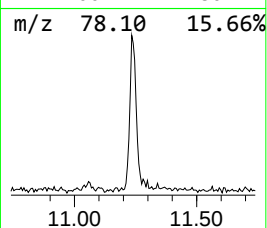
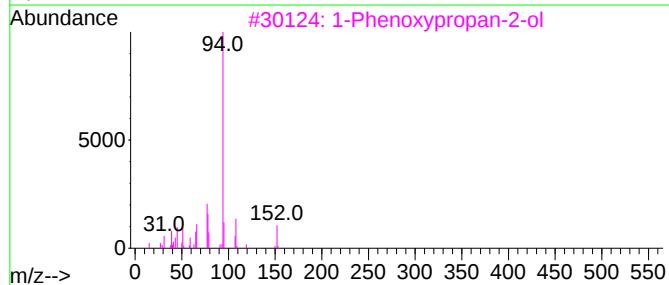
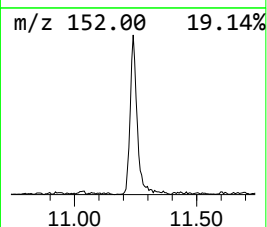
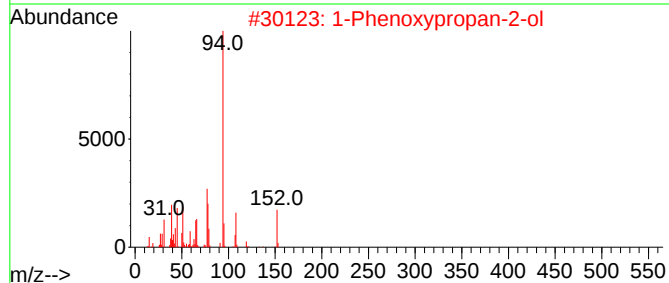
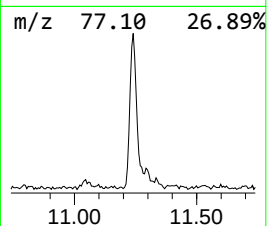
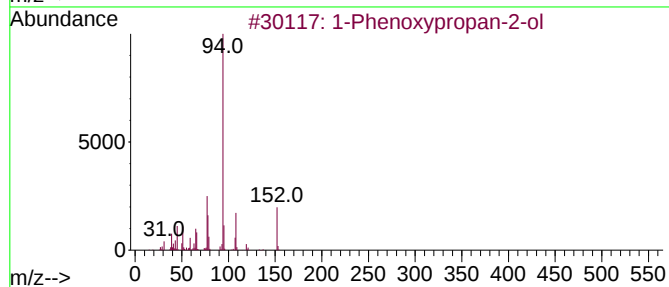
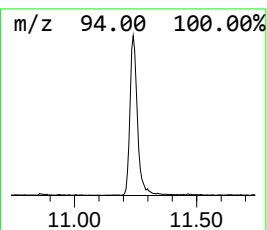
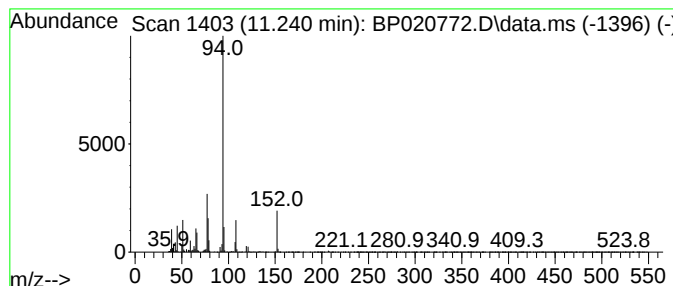
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.240	2.58 ng/ul	229062	Naphthalene-d8	10.499

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5		1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



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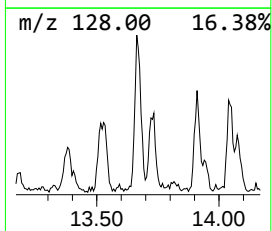
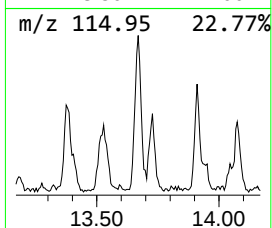
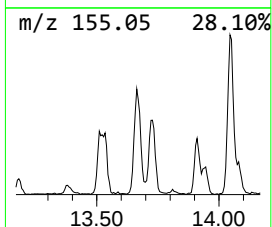
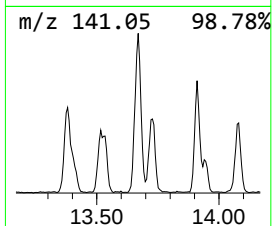
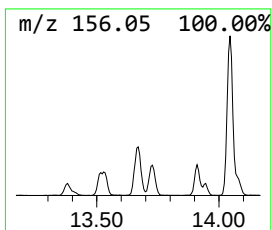
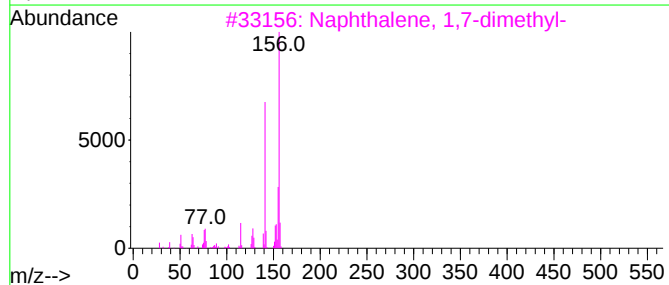
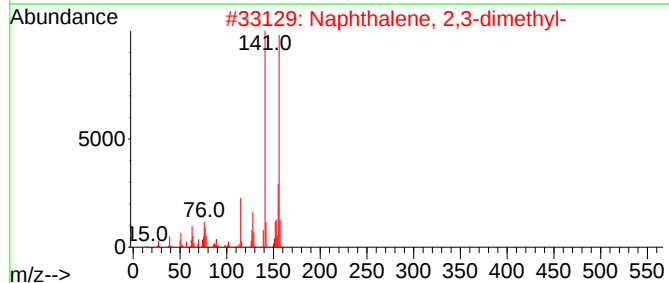
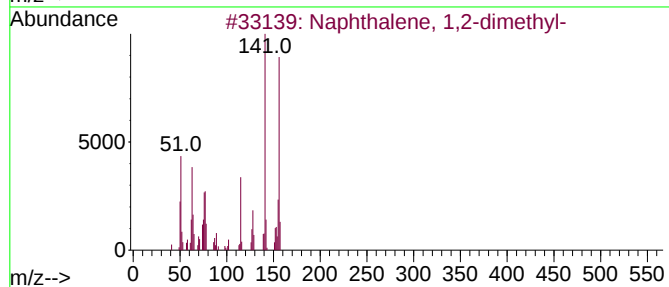
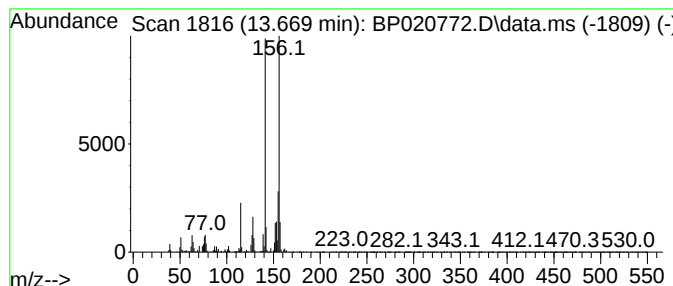
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 1,2-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.669	2.19 ng/ul	273990	Acenaphthene-d10	14.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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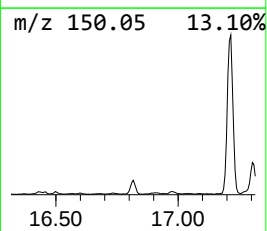
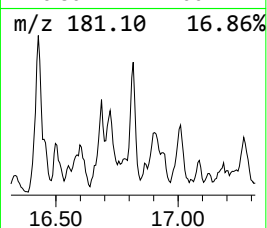
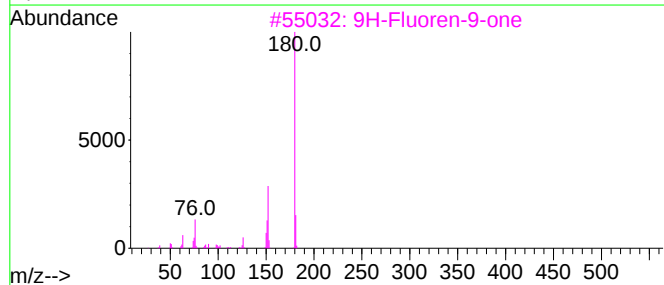
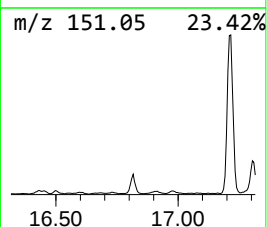
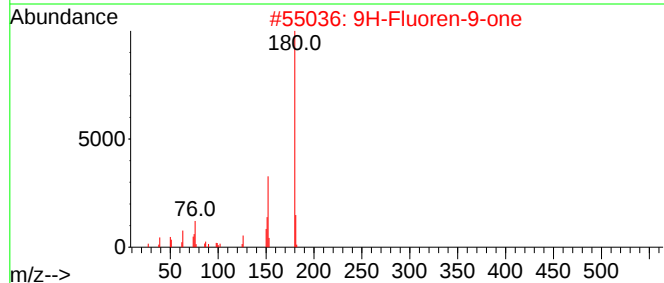
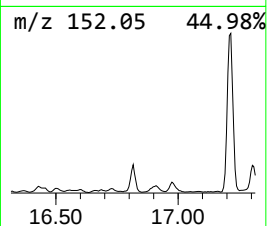
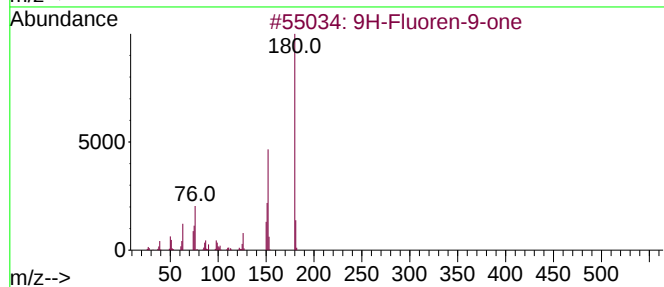
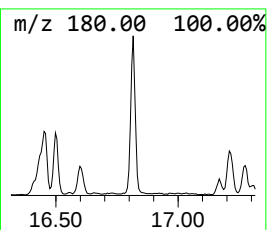
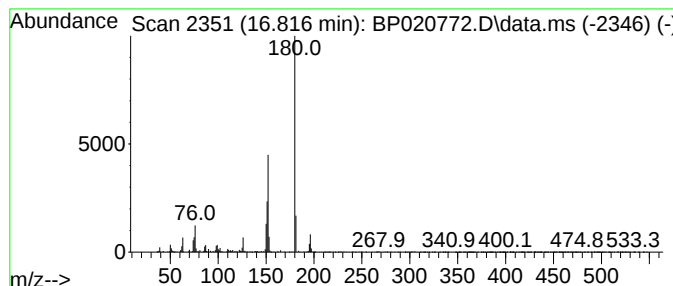
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 9H-Fluoren-9-one Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.816	2.76 ng/ul	377917	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one			180	C13H8O	000486-25-9	96
2	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
3	9H-Fluoren-9-one			180	C13H8O	000486-25-9	93
4	9,10-Phenanthredione			208	C14H8O2	000084-11-7	90
5	9,10-Phenanthredione			208	C14H8O2	000084-11-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

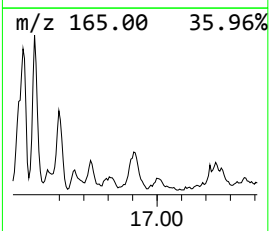
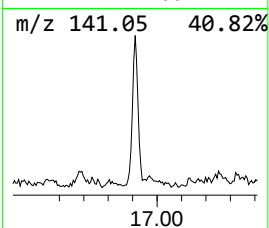
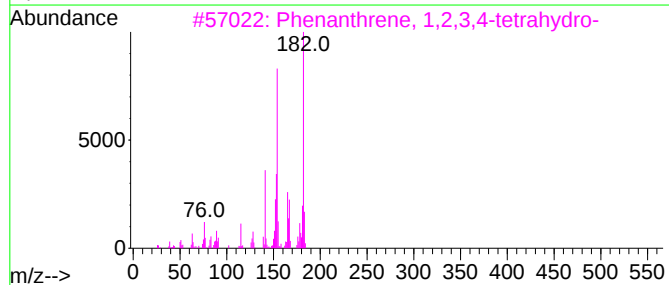
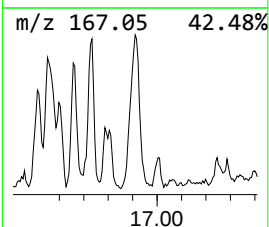
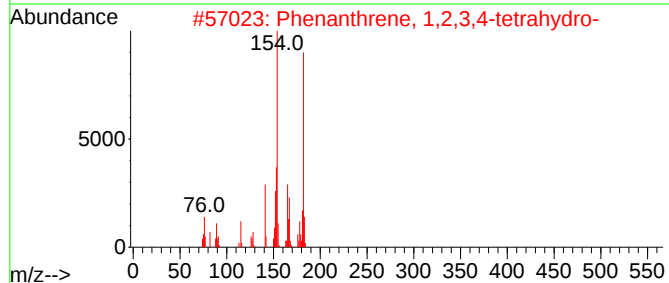
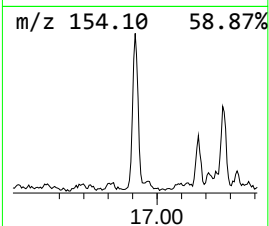
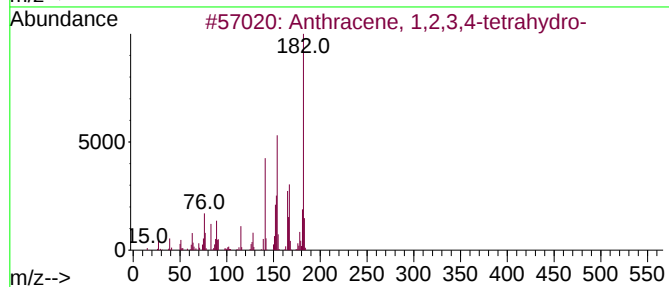
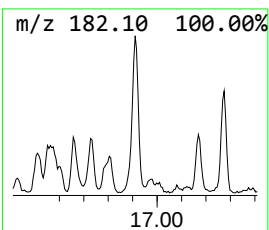
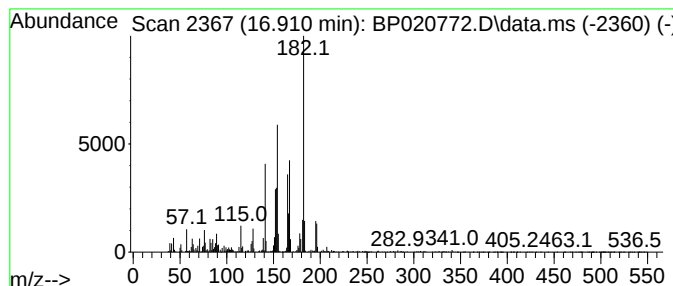
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 1,2,3,4-tetrahy... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.910	3.77 ng/ul	516577	Phenanthrene-d10		17.169	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Anthracene, 1,2,3,4-tetrahydro-		182	C14H14	002141-42-6	93
2	Phenanthrene, 1,2,3,4-tetrahydro-		182	C14H14	001013-08-7	83
3	Phenanthrene, 1,2,3,4-tetrahydro-		182	C14H14	001013-08-7	64
4	Anthracene, 1,2,3,4-tetrahydro-		182	C14H14	002141-42-6	58
5	3,3'-Dimethylbiphenyl		182	C14H14	000612-75-9	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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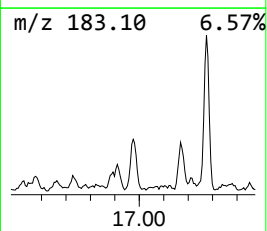
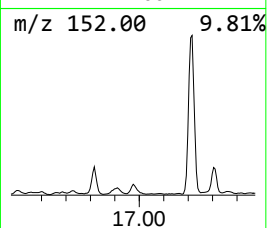
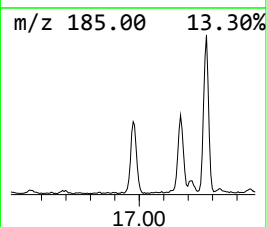
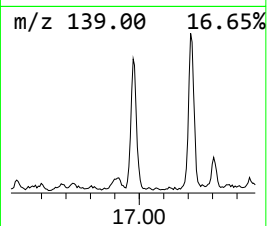
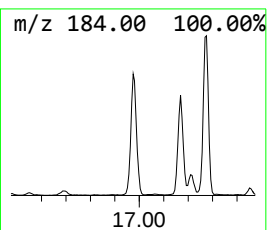
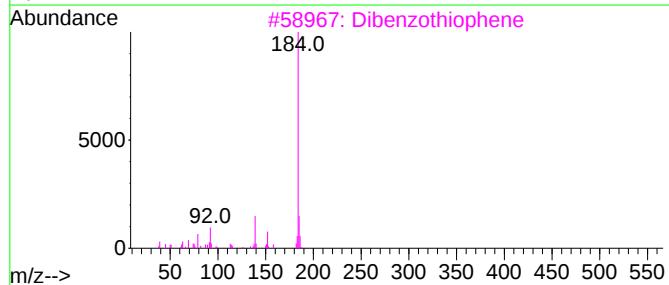
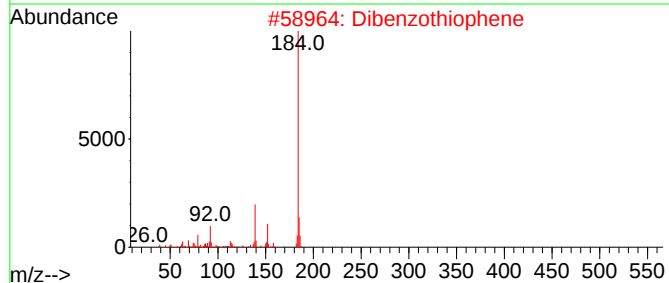
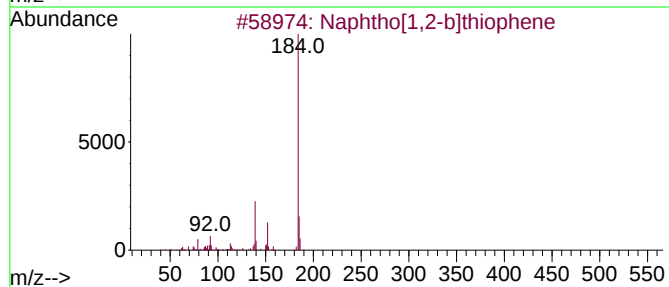
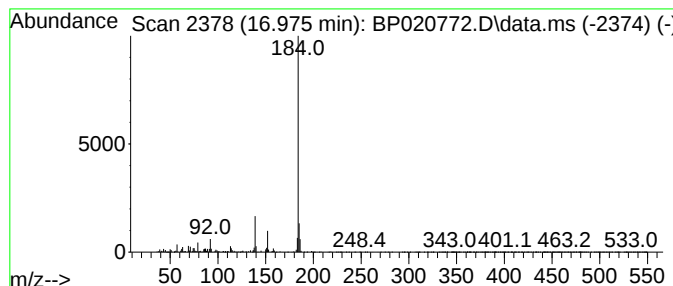
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphtho[1,2-b]thiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.975	4.20 ng/ul	574371	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	95
2			Dibenzothiophene	184	C12H8S	000132-65-0	95
3			Dibenzothiophene	184	C12H8S	000132-65-0	95
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	94
5			Dibenzothiophene	184	C12H8S	000132-65-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
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Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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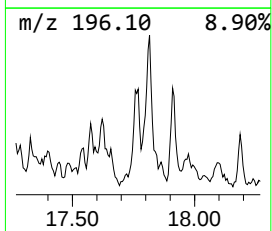
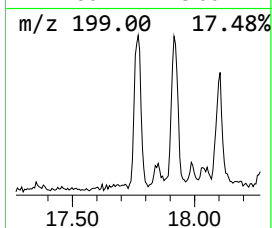
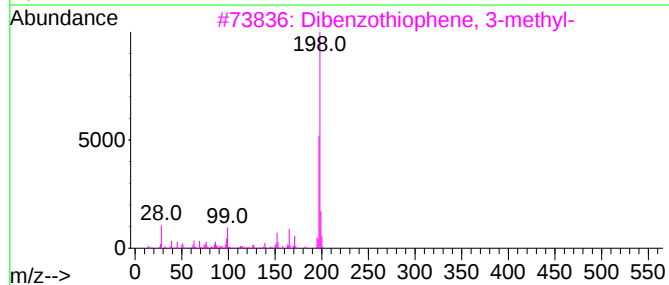
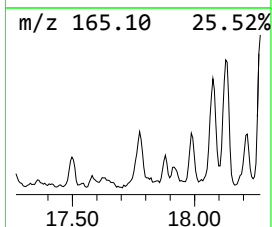
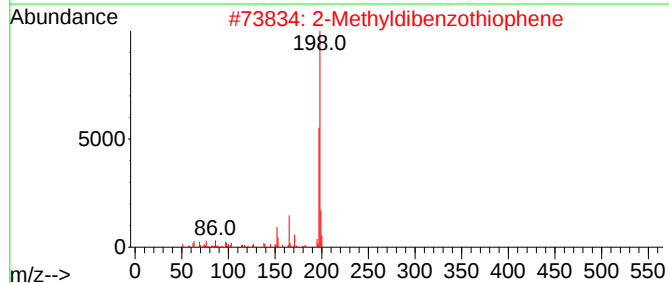
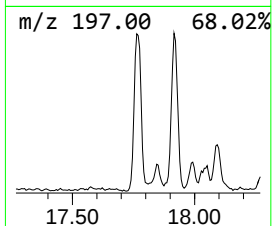
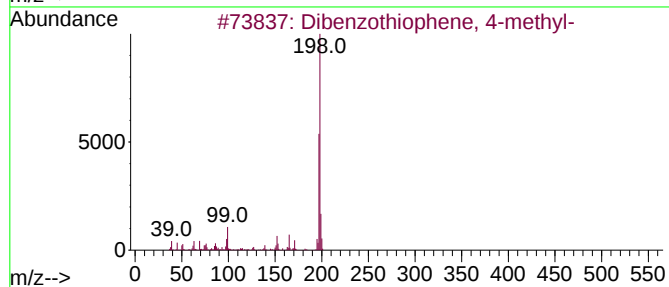
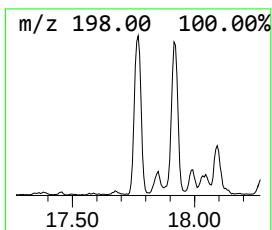
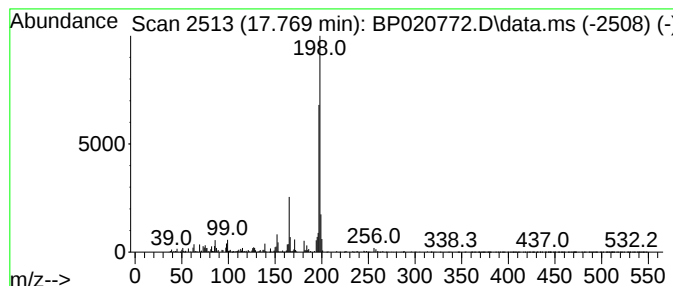
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Dibenzothiophene, 4-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.769	3.03 ng/ul	415443	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	87
2			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	87
3			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	87
4			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	86
5			4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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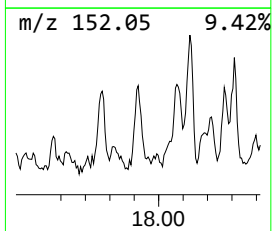
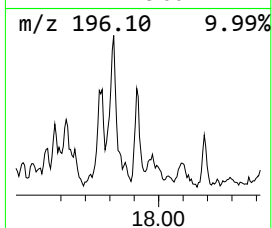
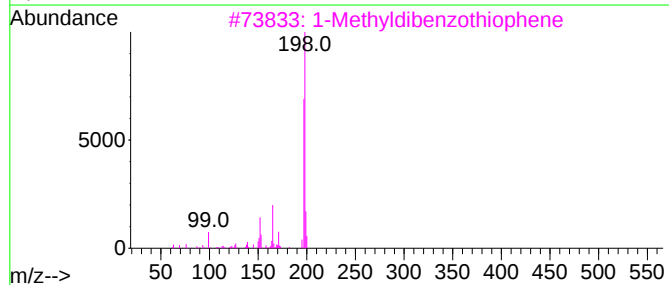
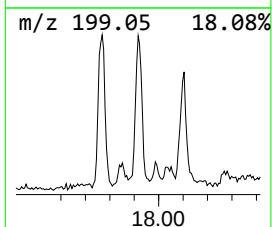
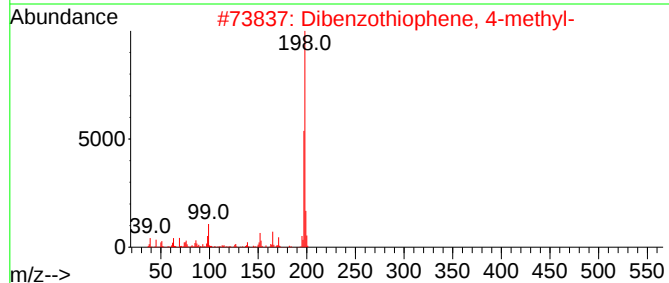
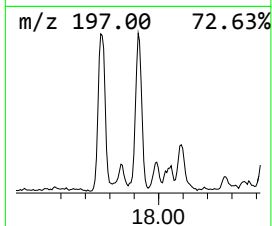
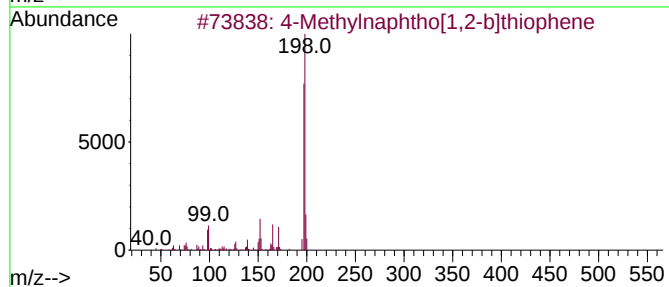
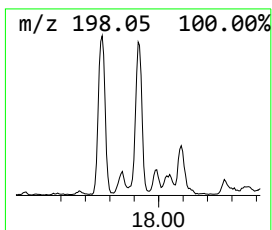
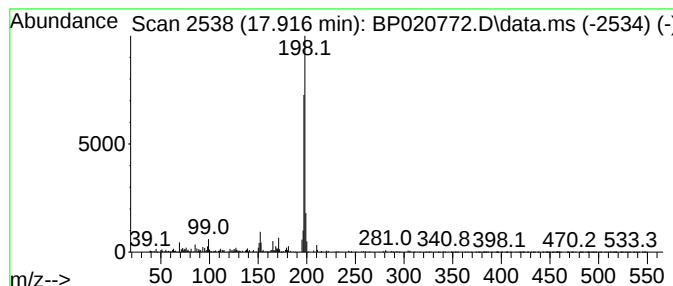
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4-Methylnaphtho[1,2-b]thiop... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.916	2.16 ng/ul	295082	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	93
2			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	90
3			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	90
4			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	90
5			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
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Misc :
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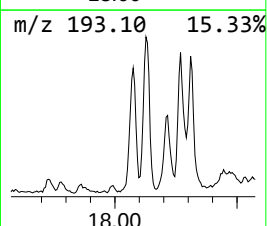
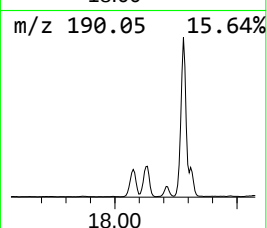
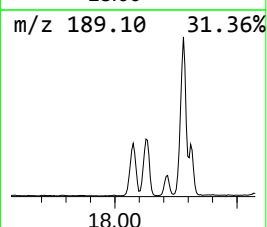
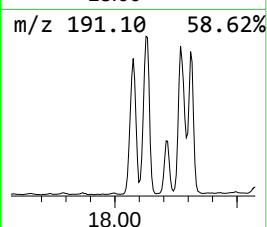
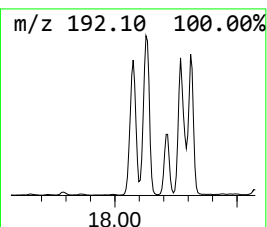
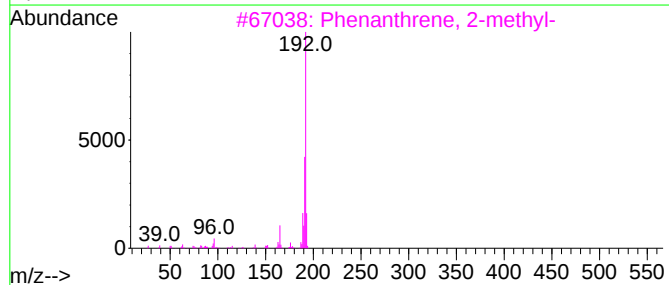
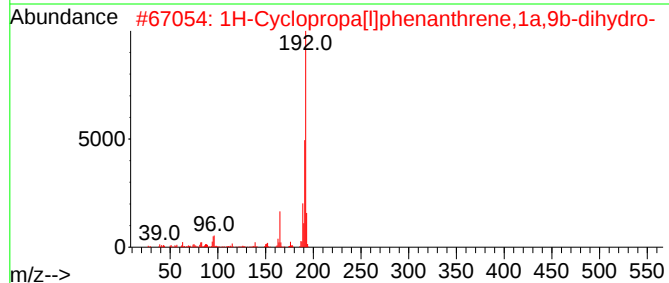
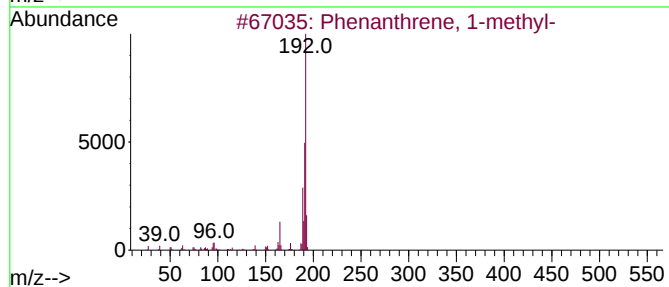
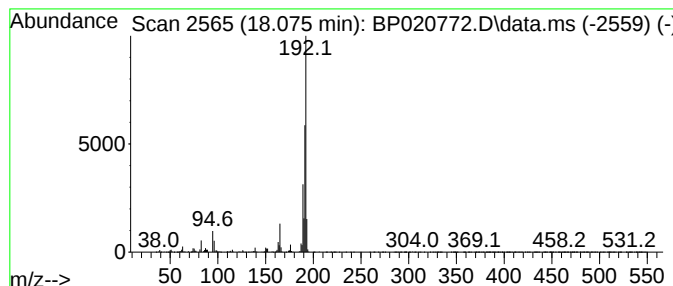
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.075	10.16 ng/ul	1390510	Phenanthrene-d10	17.169

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
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ClientSampleId :
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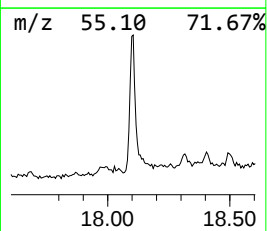
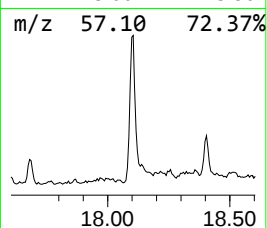
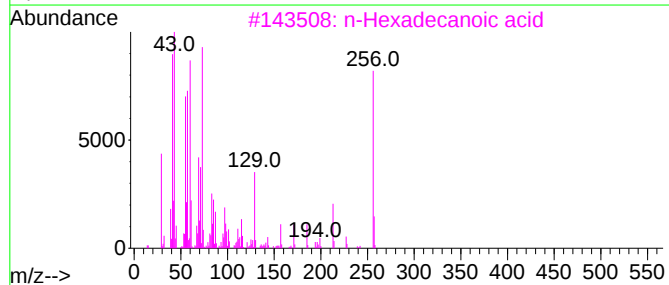
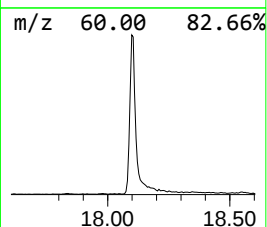
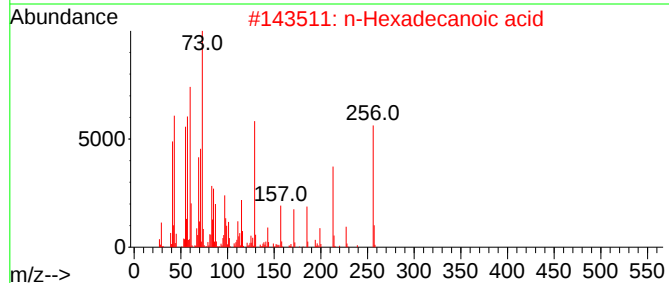
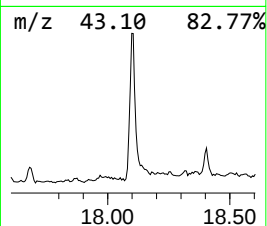
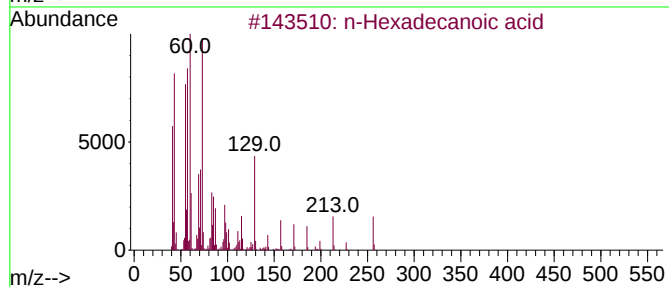
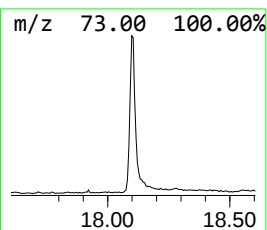
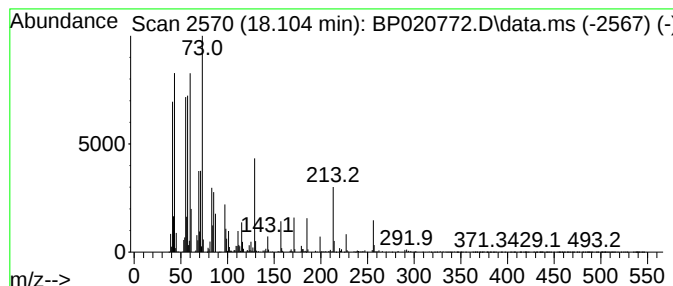
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	8.85 ng/ul	1211050	Phenanthrene-d10	17.169

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	98
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
5		Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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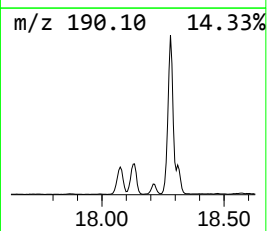
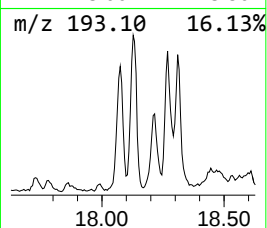
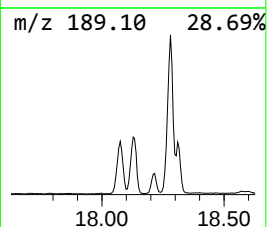
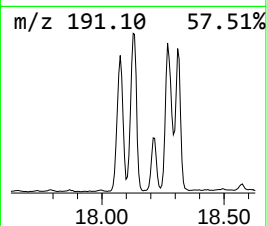
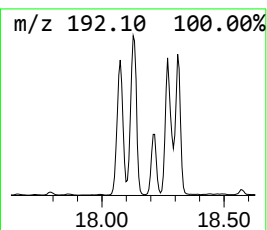
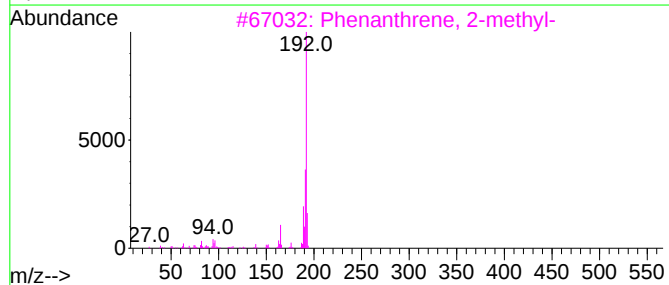
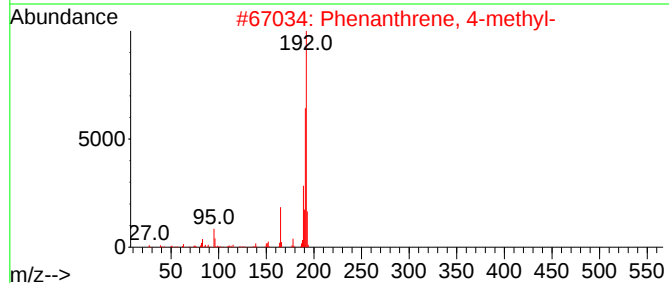
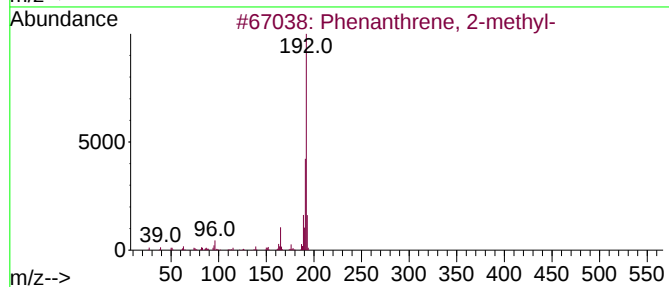
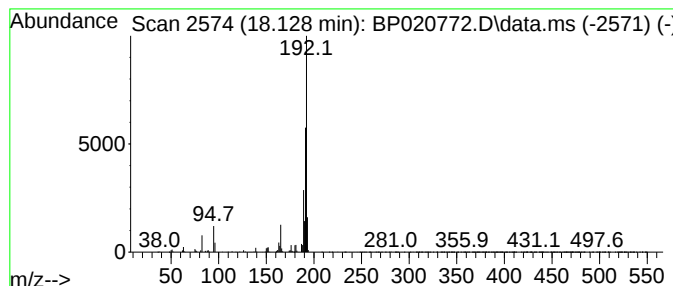
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenanthrene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	13.77 ng/ul	1885620	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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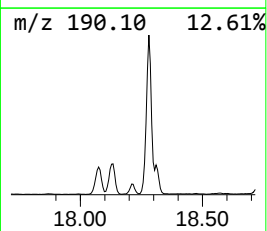
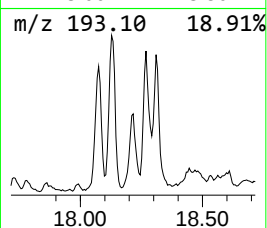
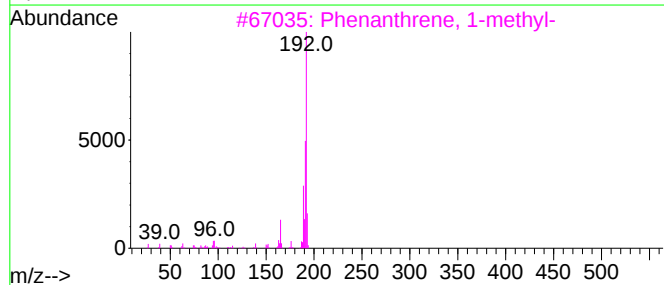
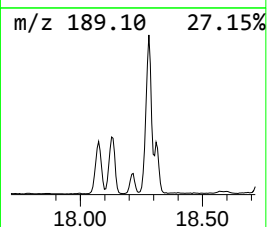
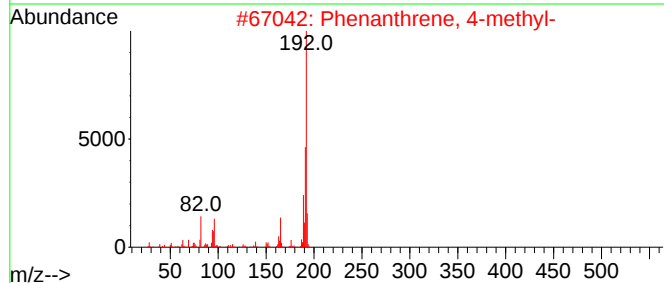
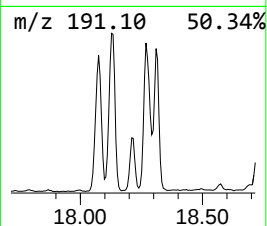
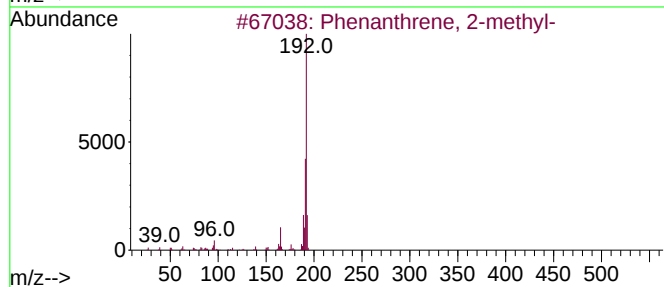
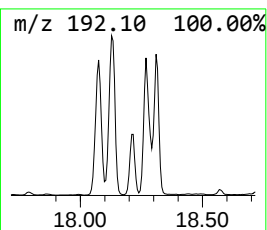
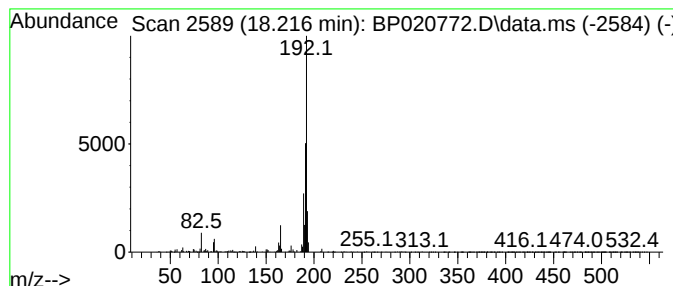
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenanthrene, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	3.95 ng/ul	540258	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
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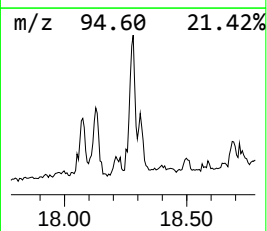
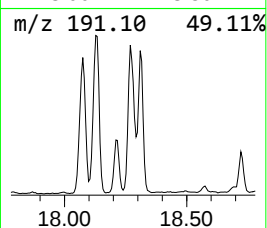
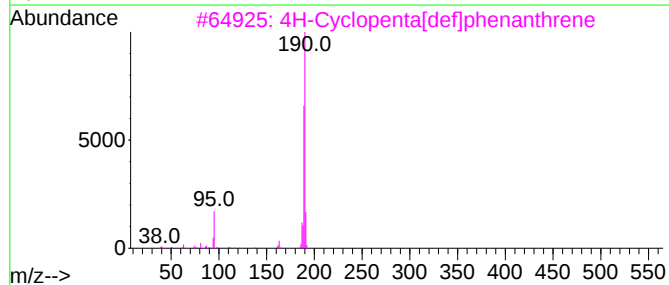
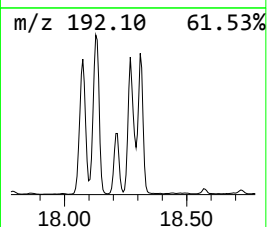
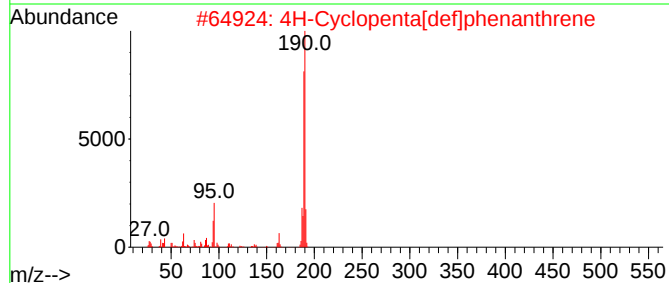
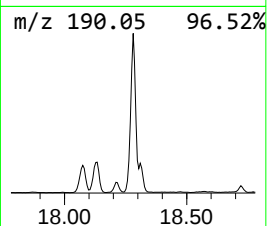
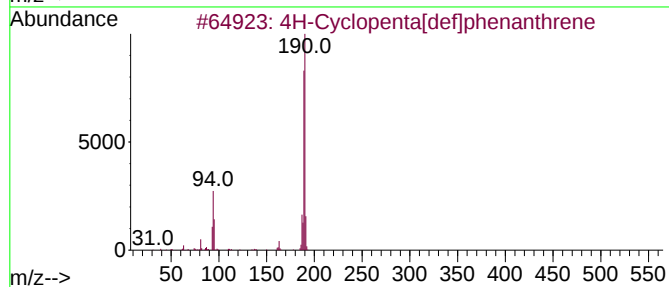
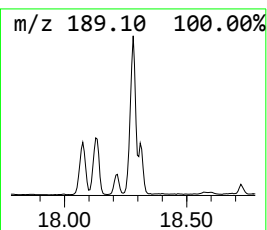
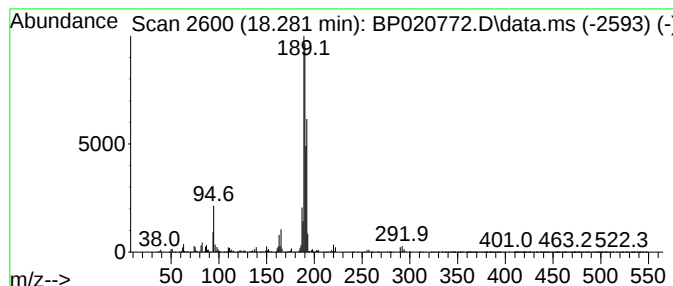
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.281	19.28 ng/ul	2640260	Phenanthrene-d10			17.169
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	70
2	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	60
3	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	43
4	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	42
5	Carbonic acid, ethyl 2-formyl-4,...		262	C10H8Cl2O4	1000331-34-4	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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Acq On : 03 Jul 2024 15:56
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Sample : P2963-13
Misc :
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Instrument :
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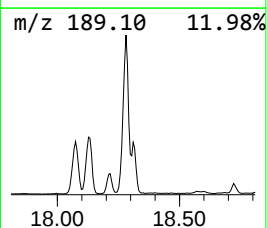
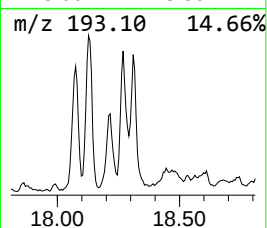
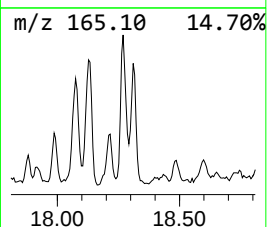
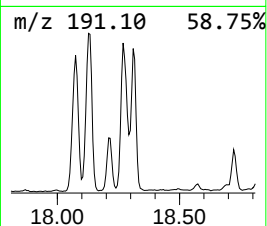
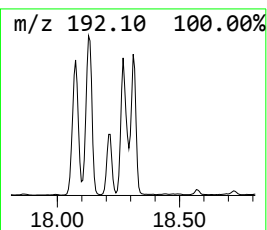
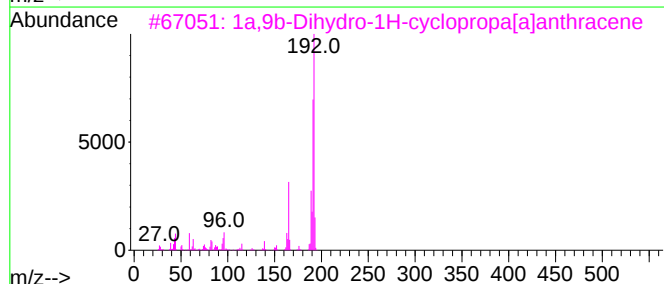
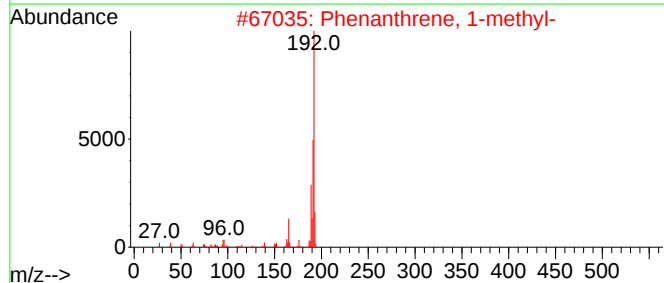
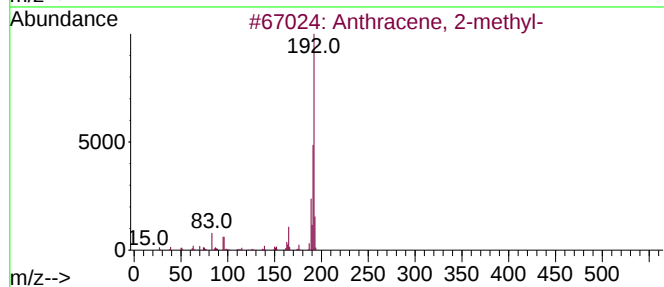
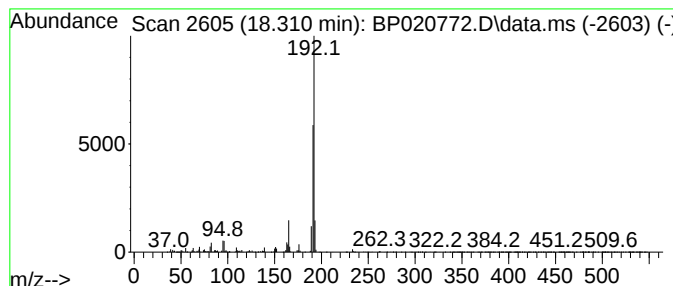
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Anthracene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	8.89 ng/ul	1217270	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	90
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
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Misc :
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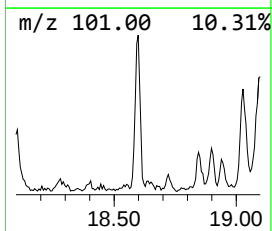
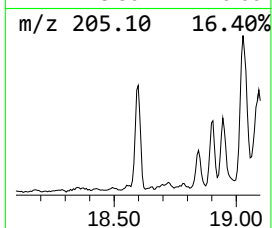
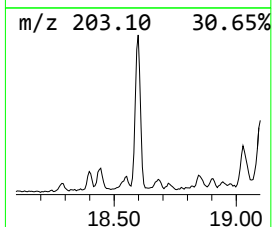
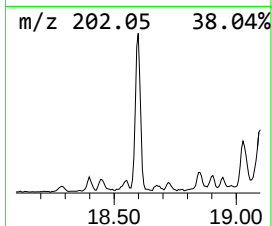
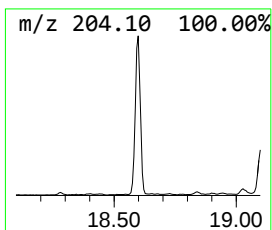
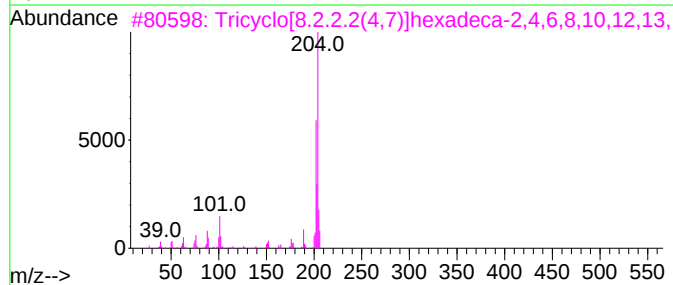
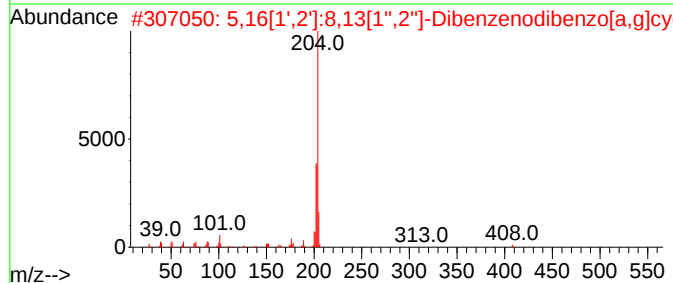
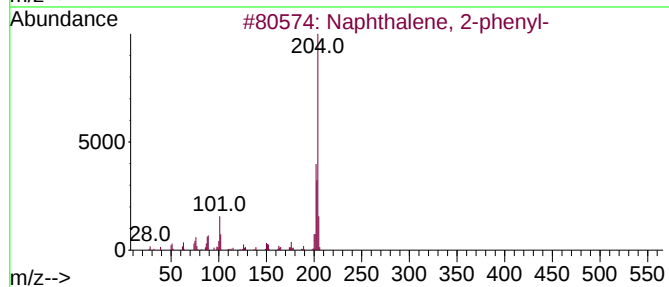
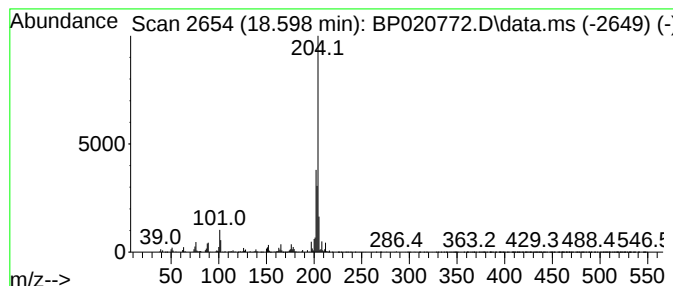
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.598	5.73 ng/ul	783976	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81
5			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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Acq On : 03 Jul 2024 15:56
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Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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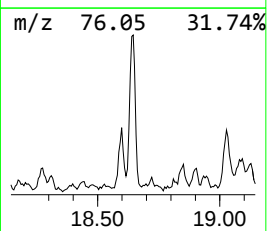
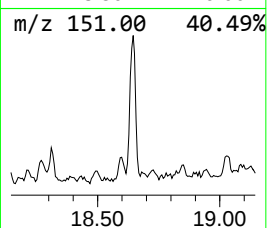
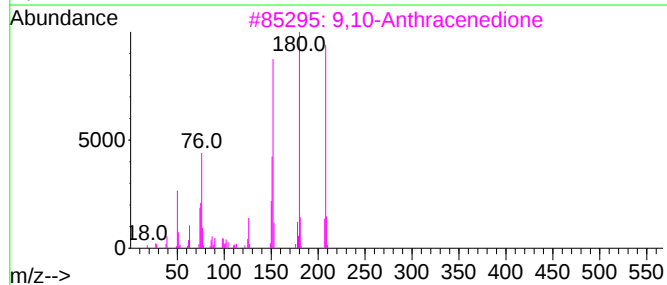
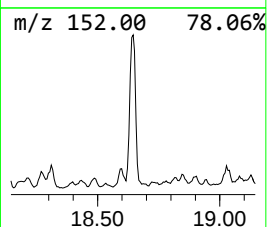
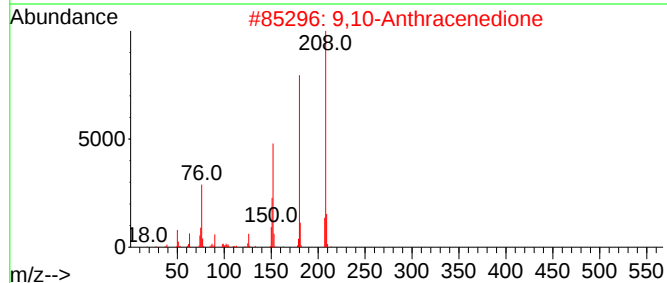
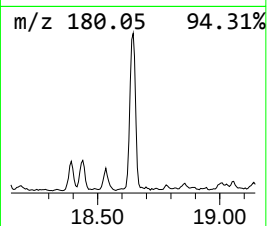
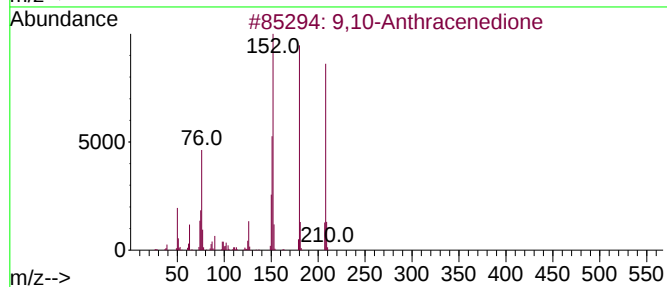
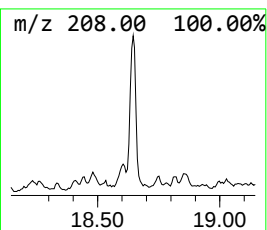
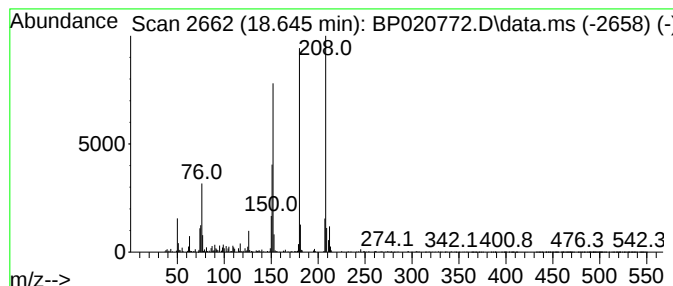
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 9,10-Anthracenedione Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.645	4.63 ng/ul	633236	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	91
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	83
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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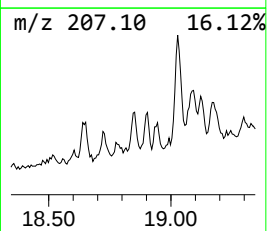
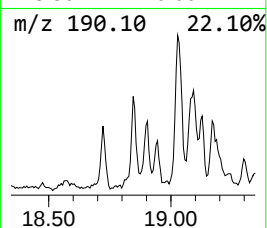
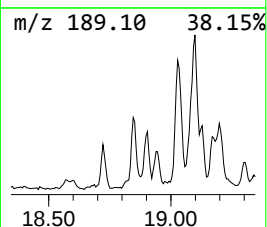
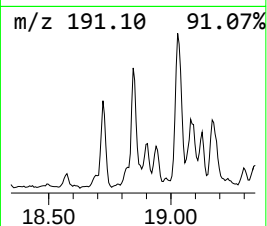
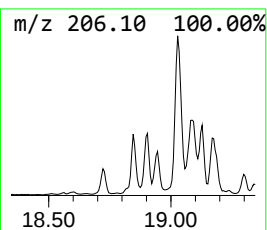
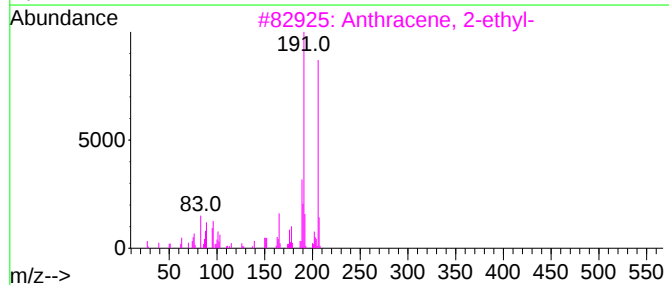
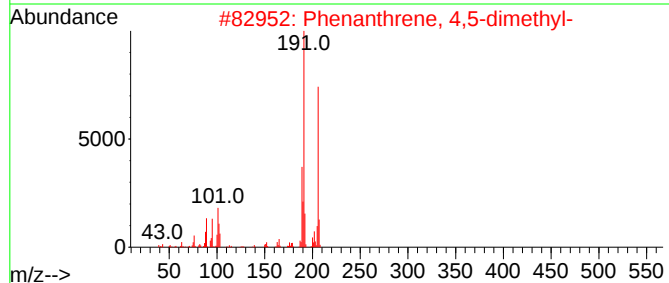
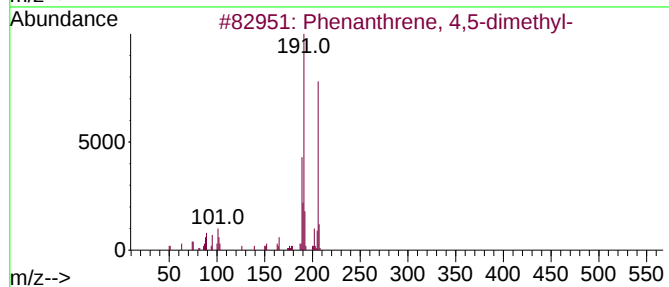
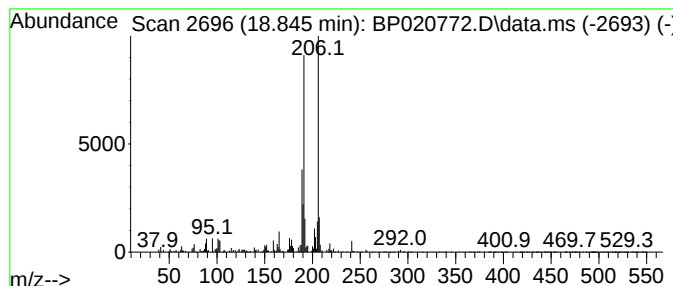
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Phenanthrene, 4,5-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.845	2.94 ng/ul	402861	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	96
2			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	94
3			Anthracene, 2-ethyl-	206	C16H14	052251-71-5	93
4			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	93
5			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
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Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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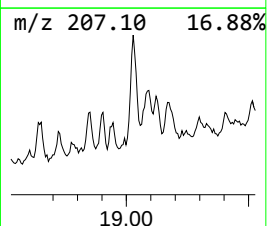
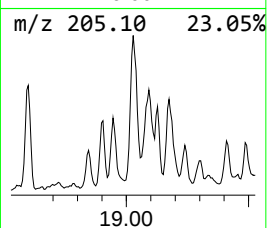
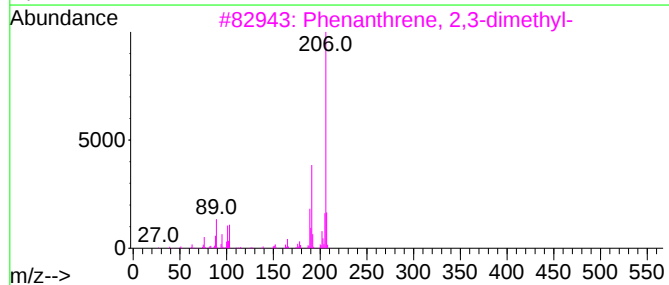
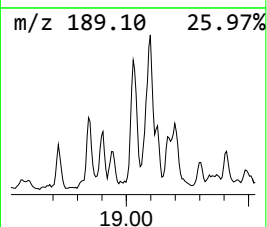
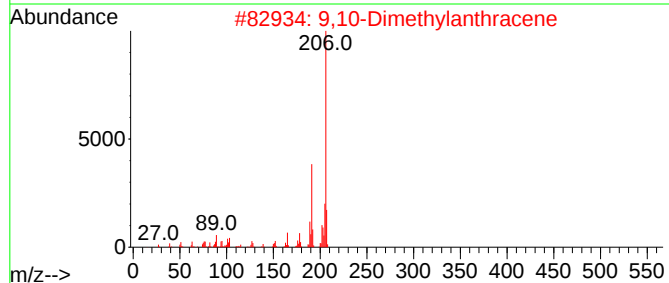
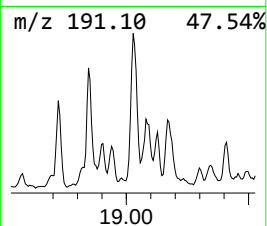
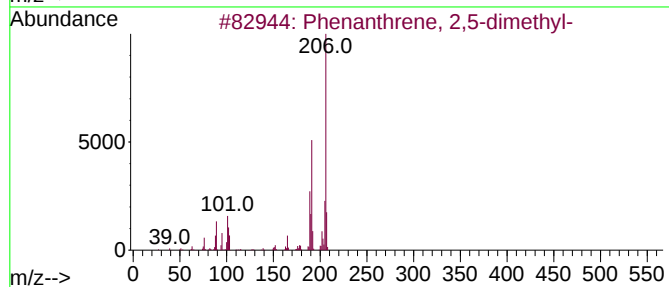
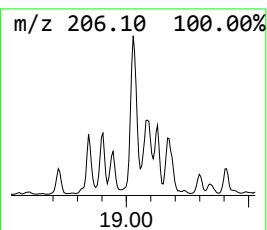
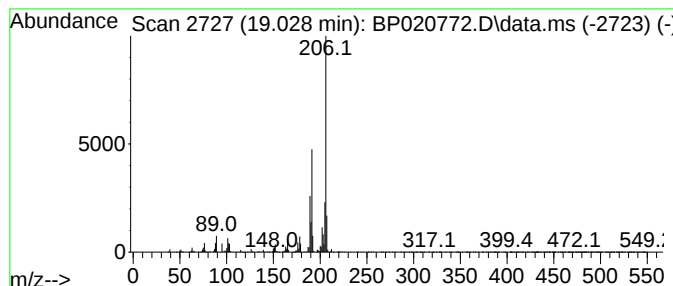
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.028	7.51 ng/ul	1028310	Phenanthrene-d10	17.169

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2		9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	91
5		9,10-Dimethylanthracene	206	C16H14	000781-43-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

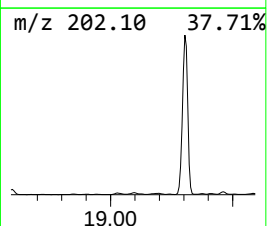
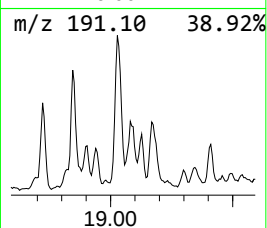
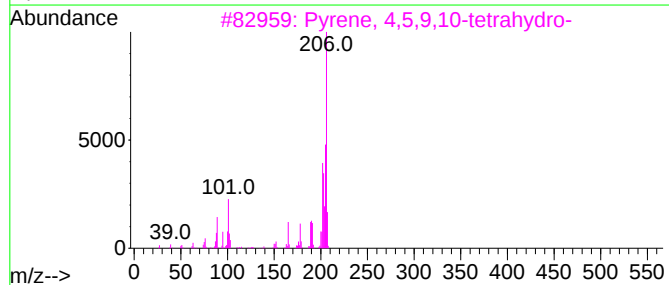
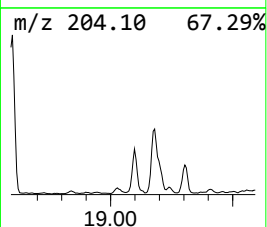
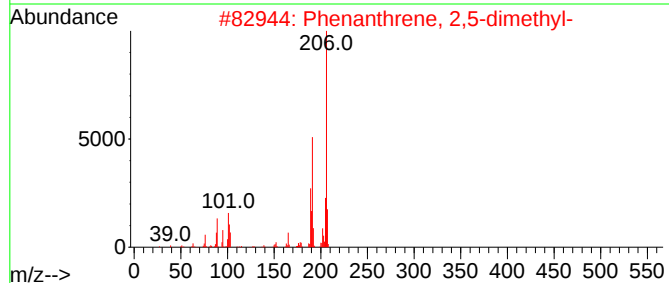
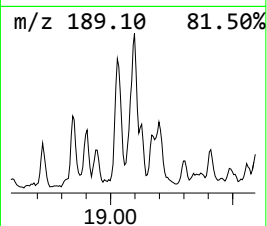
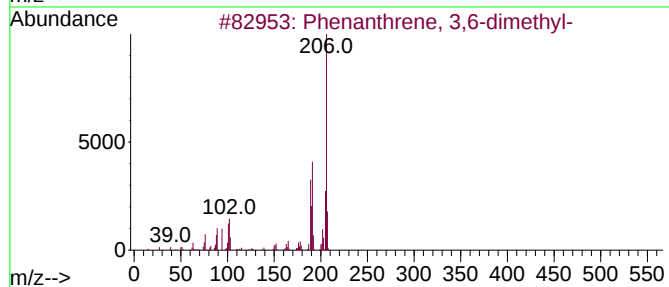
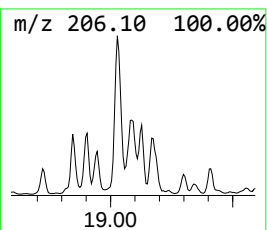
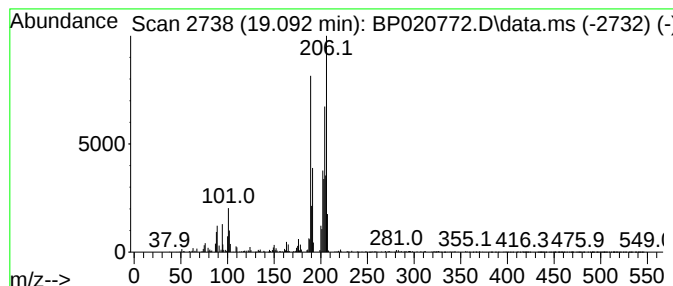
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 19 Phenanthrene, 3,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.092	4.98 ng/ul	681386	Phenanthrene-d10		17.169	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	50
2	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	45
3	Pyrene, 4,5,9,10-tetrahydro-		206	C16H14	000781-17-9	42
4	Naphthalene, 1,2-dihydro-4-phenyl-		206	C16H14	007469-40-1	38
5	1,2,4,8-Tetramethylbicyclo[6.3.0...]		204	C15H24	137235-51-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

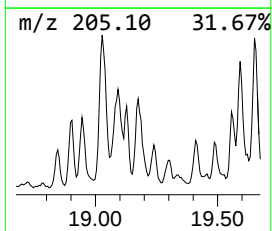
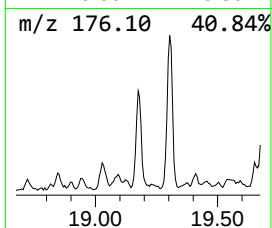
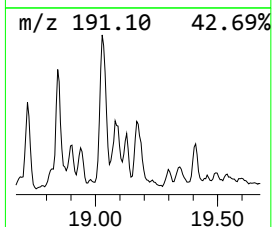
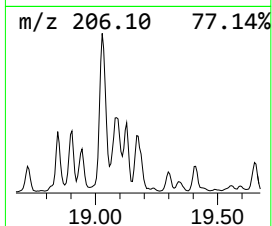
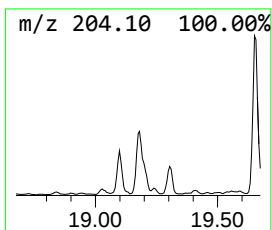
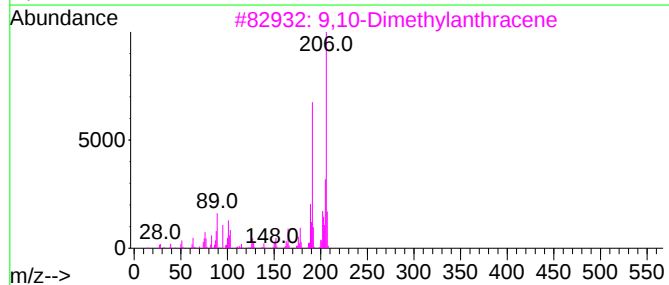
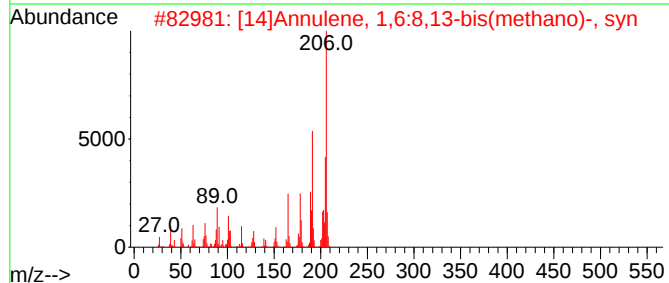
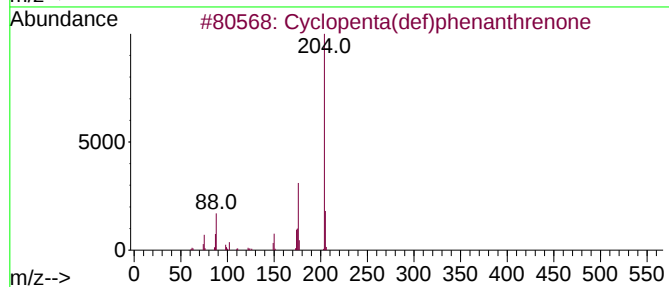
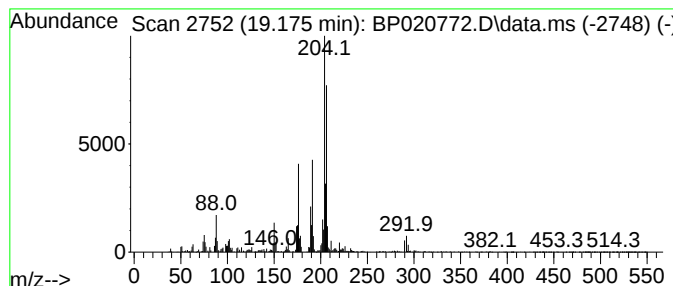
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 Cyclopenta(def)phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.175	8.26 ng/ul	1130440	Phenanthrene-d10	17.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	83
2	[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	70
3	9,10-Dimethylantracene	206	C16H14	000781-43-1	55
4	9,10-Dimethylantracene	206	C16H14	000781-43-1	50
5	Benzene, (2-methylene-1-phenylcy...	206	C16H14	025152-47-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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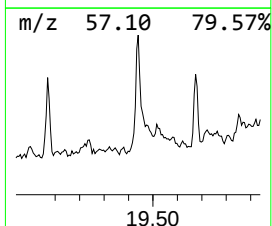
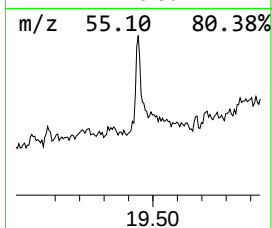
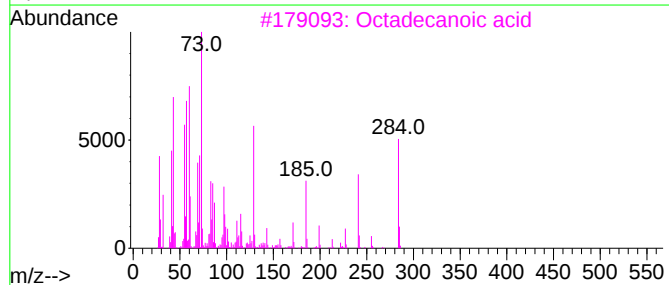
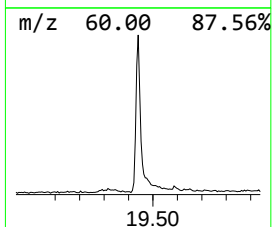
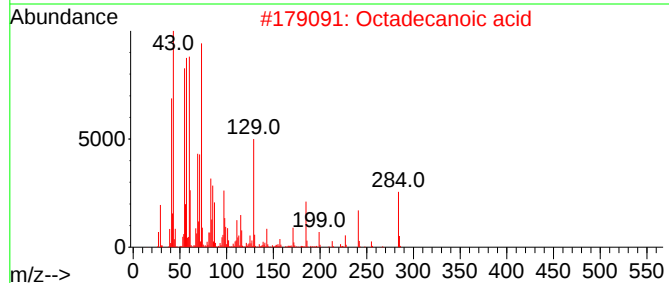
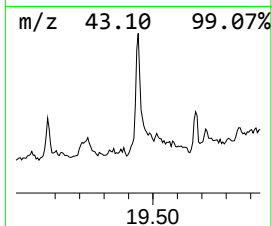
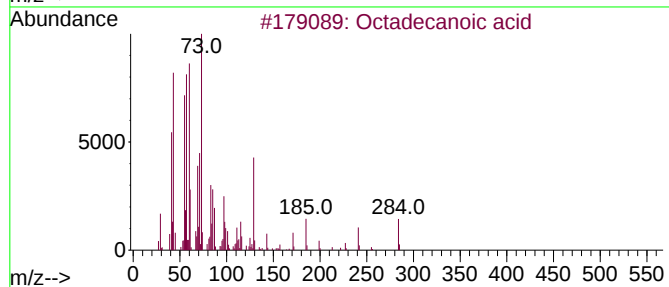
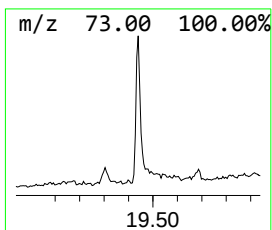
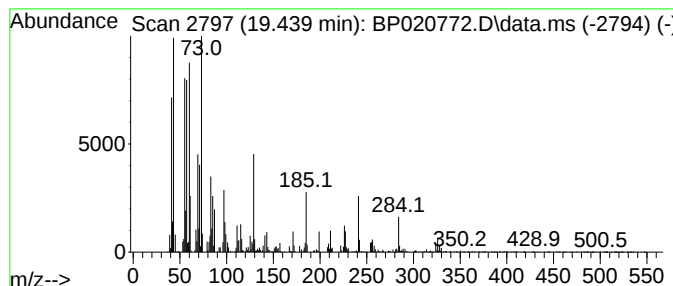
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Octadecanoic acid Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.439	2.35 ng/ul	773048	Chrysene-d12	21.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	96
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
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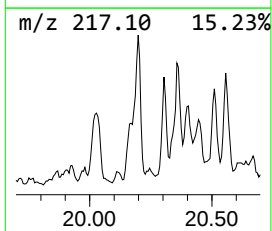
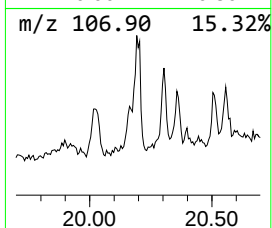
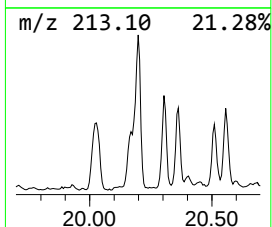
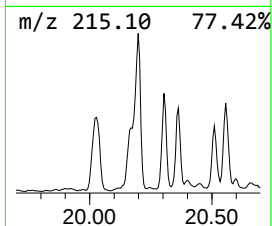
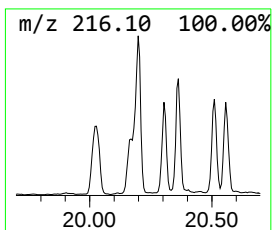
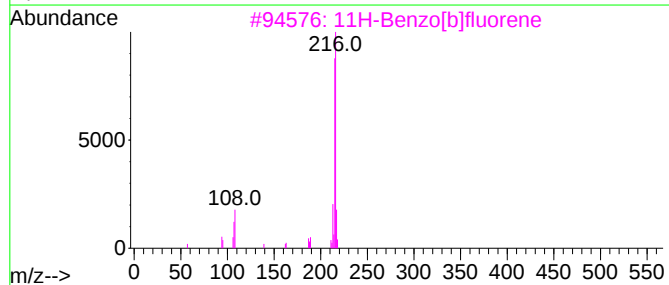
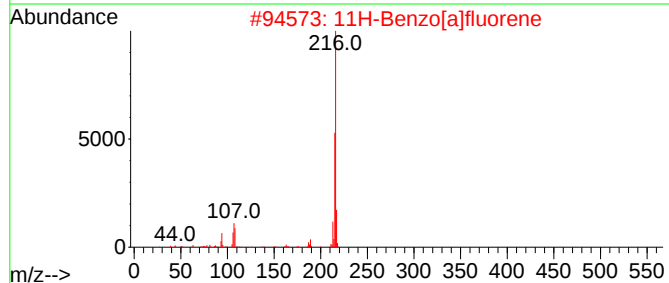
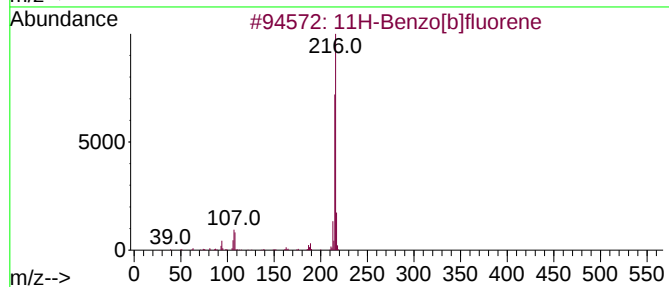
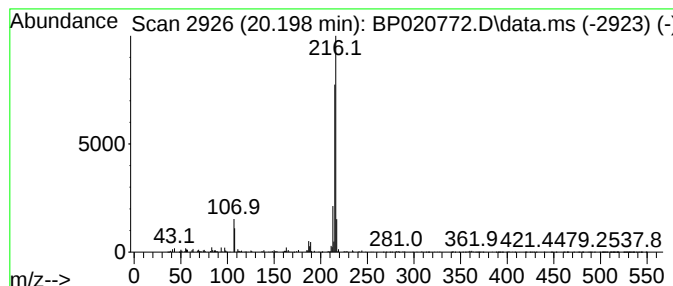
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 11H-Benzo[b]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.198	3.30 ng/ul	1084360	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene			216	C17H12	000243-17-4	91
2	11H-Benzo[a]fluorene			216	C17H12	000238-84-6	90
3	11H-Benzo[b]fluorene			216	C17H12	000243-17-4	87
4	11H-Benzo[b]fluorene			216	C17H12	000243-17-4	87
5	11H-Benzo[a]fluorene			216	C17H12	000238-84-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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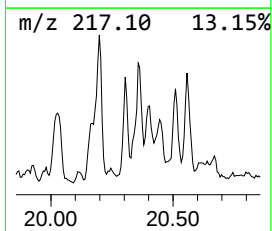
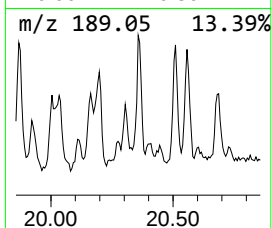
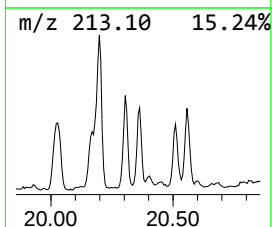
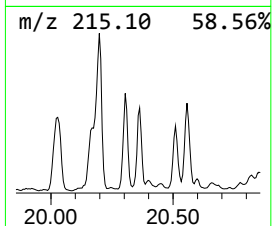
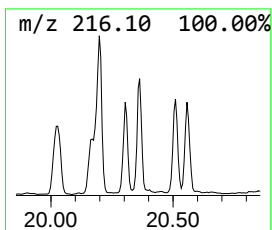
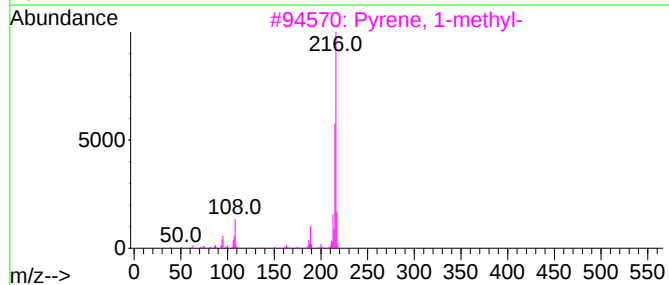
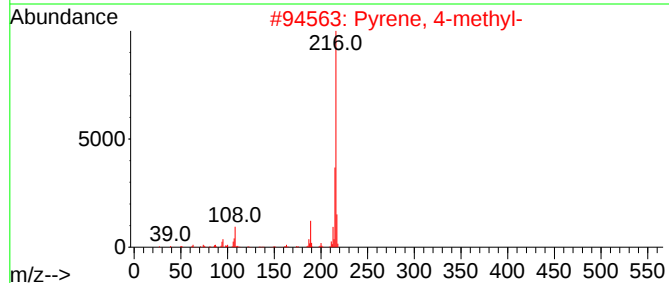
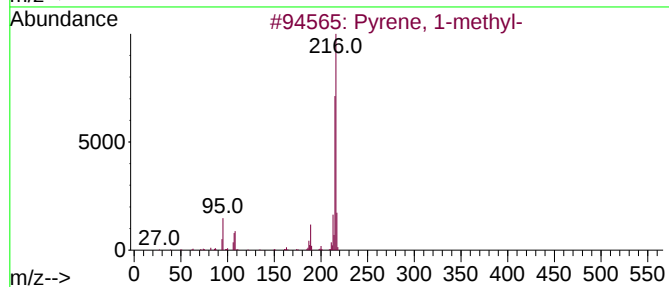
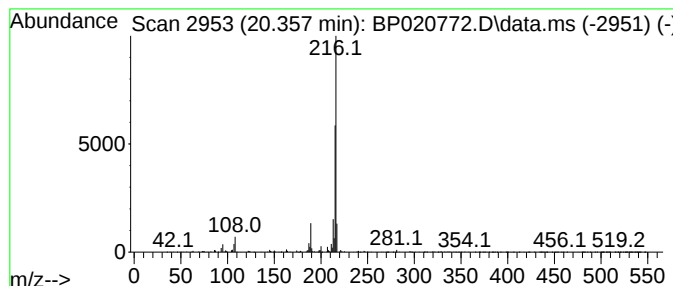
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Pyrene, 1-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.357	2.07 ng/ul	679992	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2			Pyrene, 4-methyl-	216	C17H12	003353-12-6	95
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CR7

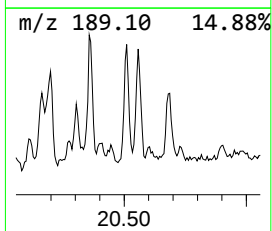
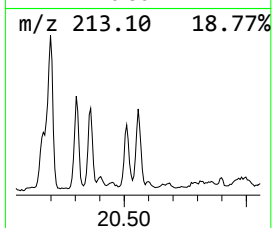
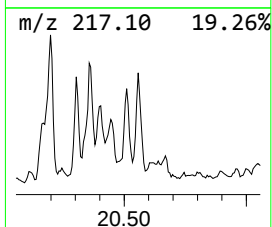
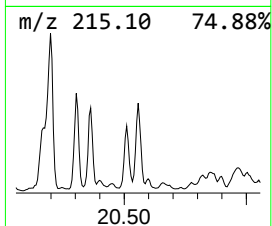
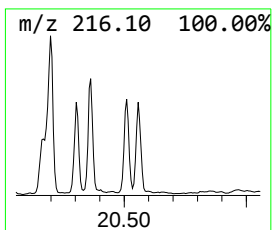
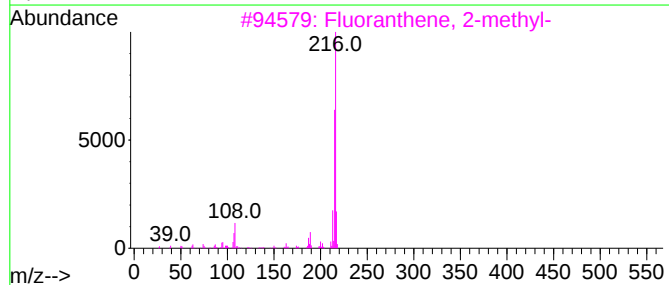
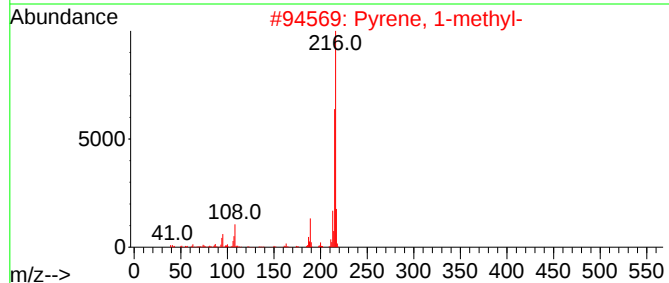
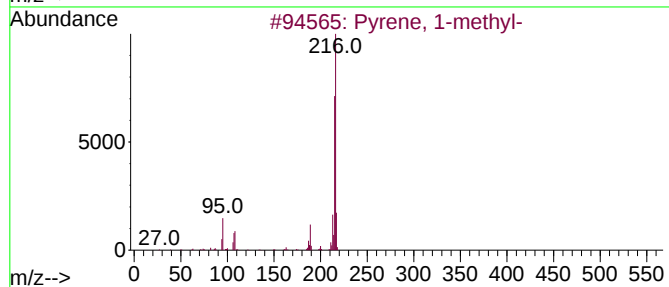
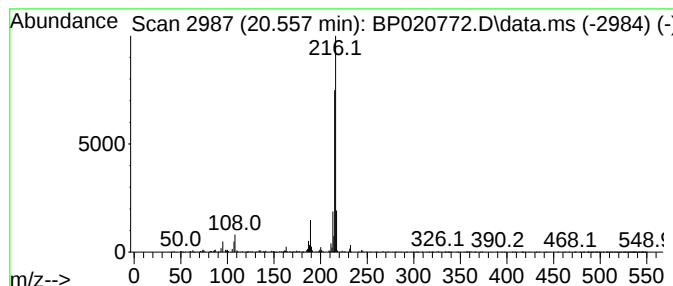
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Fluoranthene, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.557	2.27 ng/ul	746535	Chrysene-d12	21.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
5		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
Data File : BP020772.D
Acq On : 03 Jul 2024 15:56
Operator : MA/JU
Sample : P2963-13
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CR7

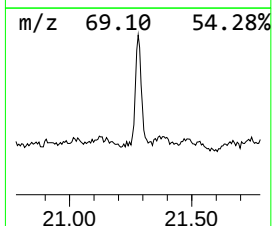
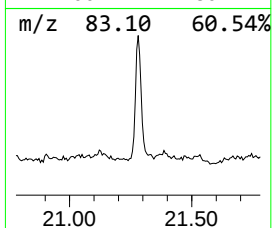
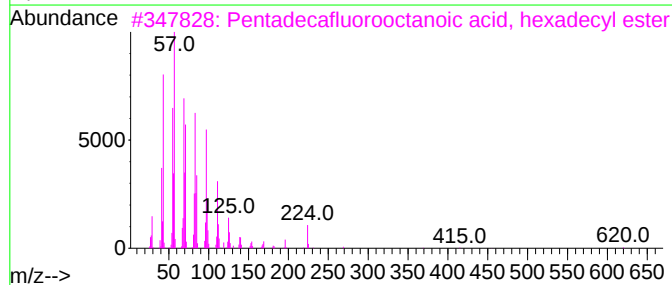
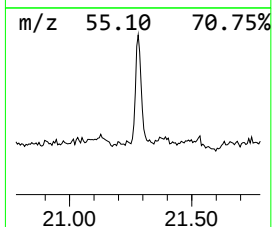
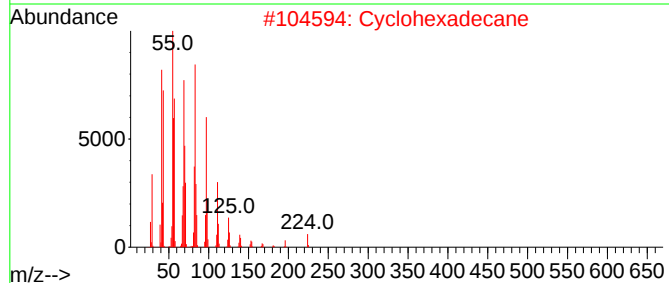
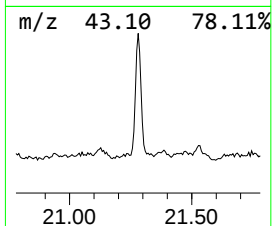
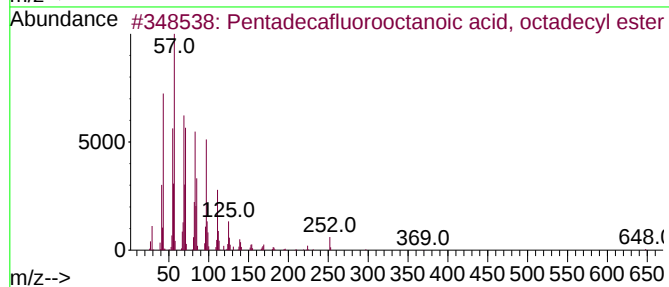
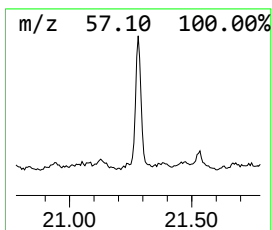
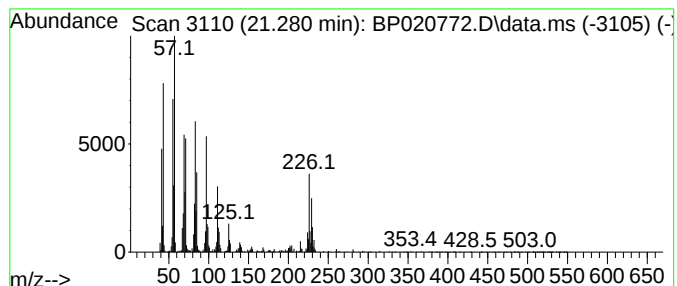
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Pentadecafluorooctanoic aci... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.280	6.72 ng/ul	2207710	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2			Cyclohexadecane	224	C16H32	000295-65-8	90
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	89
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	78
5			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070324\
 Data File : BP020772.D
 Acq On : 03 Jul 2024 15:56
 Operator : MA/JU
 Sample : P2963-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CR7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.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
Silane, tetrame...	4.899	3.5	ng/ul	212468	1	7.740	1232740	20.0
1-Phenoxypropan...	11.240	2.6	ng/ul	229062	2	10.499	1775700	20.0
Naphthalene, 1,...	13.669	2.2	ng/ul	273990	3	14.351	2507610	20.0
9H-Fluoren-9-one	16.816	2.8	ng/ul	377917	4	17.169	2738150	20.0
Anthracene, 1,2...	16.910	3.8	ng/ul	516577	4	17.169	2738150	20.0
Naphtho[1,2-b]t...	16.975	4.2	ng/ul	574371	4	17.169	2738150	20.0
Dibenzothiophen...	17.769	3.0	ng/ul	415443	4	17.169	2738150	20.0
4-Methylnaphtho...	17.916	2.2	ng/ul	295082	4	17.169	2738150	20.0
Phenanthrene, 1...	18.075	10.2	ng/ul	1390510	4	17.169	2738150	20.0
n-Hexadecanoic ...	18.104	8.8	ng/ul	1211050	4	17.169	2738150	20.0
Phenanthrene, 2...	18.128	13.8	ng/ul	1885620	4	17.169	2738150	20.0
Phenanthrene, 4...	18.216	4.0	ng/ul	540258	4	17.169	2738150	20.0
4H-Cyclopenta[d...	18.281	19.3	ng/ul	2640260	4	17.169	2738150	20.0
Anthracene, 2-m...	18.310	8.9	ng/ul	1217270	4	17.169	2738150	20.0
Naphthalene, 2-...	18.598	5.7	ng/ul	783976	4	17.169	2738150	20.0
9,10-Anthracene...	18.645	4.6	ng/ul	633236	4	17.169	2738150	20.0
Phenanthrene, 4...	18.845	2.9	ng/ul	402861	4	17.169	2738150	20.0
Phenanthrene, 2...	19.028	7.5	ng/ul	1028310	4	17.169	2738150	20.0
Phenanthrene, 3...	19.092	5.0	ng/ul	681386	4	17.169	2738150	20.0
Cyclopenta(def)...	19.175	8.3	ng/ul	1130440	4	17.169	2738150	20.0
Octadecanoic acid	19.439	2.4	ng/ul	773048	5	21.622	6567650	20.0
11H-Benzo[b]flu...	20.198	3.3	ng/ul	1084360	5	21.622	6567650	20.0
Pyrene, 1-methyl-	20.357	2.1	ng/ul	679992	5	21.622	6567650	20.0
Fluoranthene, 2...	20.557	2.3	ng/ul	746535	5	21.622	6567650	20.0
Pentadecafluoro...	21.280	6.7	ng/ul	2207710	5	21.622	6567650	20.0