

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
 Data File : BP022450.D
 Acq On : 17 Oct 2024 18:42
 Operator : RC/JU
 Sample : P4264-02MEDL2 20X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCYN6MEDL2

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022450.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.623	105	108	119	rBV	9517	17267	0.78%	0.084%
2	3.928	156	160	167	rVB2	10061	12806	0.58%	0.062%
3	6.816	646	651	654	rVV2	8998	15434	0.70%	0.075%
4	6.858	654	658	672	rVB3	15254	35473	1.61%	0.172%
5	7.011	678	684	691	rVB	14365	24110	1.09%	0.117%
6	7.211	711	718	726	rBV	18243	32683	1.48%	0.158%
7	7.663	785	795	805	rBV	338474	552256	25.07%	2.674%
8	8.034	852	858	866	rBV	14743	25521	1.16%	0.124%
9	8.193	878	885	893	rVB2	6687	13736	0.62%	0.067%
10	8.375	911	916	923	rBV	18882	34796	1.58%	0.168%
11	8.810	985	990	997	rVV	16156	31502	1.43%	0.153%
12	9.122	1037	1043	1049	rVB5	4818	10378	0.47%	0.050%
13	9.222	1054	1060	1065	rVV2	10372	15604	0.71%	0.076%
14	9.522	1105	1111	1119	rBV3	12718	24665	1.12%	0.119%
15	9.828	1157	1163	1172	rVB3	7144	14626	0.66%	0.071%
16	10.069	1198	1204	1211	rBV2	18528	38046	1.73%	0.184%
17	10.422	1256	1264	1268	rBV	375243	674026	30.60%	3.264%
18	10.475	1268	1273	1283	rVV	962825	1641565	74.53%	7.949%
19	10.587	1285	1292	1303	rVV3	48631	107030	4.86%	0.518%
20	11.699	1475	1481	1485	rBV2	7609	11650	0.53%	0.056%
21	12.087	1539	1547	1557	rVV	364110	562059	25.52%	2.722%
22	12.199	1561	1566	1571	rVV3	7139	13892	0.63%	0.067%
23	12.304	1572	1584	1595	rVB	167138	276566	12.56%	1.339%
24	13.104	1713	1720	1726	rBV	101141	146334	6.64%	0.709%
25	13.304	1748	1754	1764	rVB	31298	60016	2.72%	0.291%
26	13.440	1768	1777	1788	rBV4	47888	130051	5.90%	0.630%
27	13.598	1797	1804	1809	rBV2	63862	120696	5.48%	0.584%
28	13.651	1809	1813	1816	rVV	42598	66701	3.03%	0.323%
29	13.693	1816	1820	1829	rVV2	47444	94000	4.27%	0.455%
30	13.840	1837	1845	1848	rVV	28186	46671	2.12%	0.226%
31	13.869	1848	1850	1855	rVB3	10483	11926	0.54%	0.058%
32	13.975	1862	1868	1872	rBV	48604	81164	3.68%	0.393%
33	14.004	1872	1873	1881	rVB3	21623	27447	1.25%	0.133%
34	14.287	1910	1921	1927	rBV	618231	1038096	47.13%	5.027%
35	14.351	1927	1932	1941	rVB	279573	435550	19.77%	2.109%
36	14.428	1941	1945	1951	rVB	16496	25441	1.16%	0.123%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Title : SVOA CALIBRATION

37	14.498	1951	1957	1961	rBV2	10524	24994	1.13%	0.121%
38	14.598	1967	1974	1979	rVB3	11172	22302	1.01%	0.108%
39	14.693	1979	1990	1999	rBV	320078	513727	23.32%	2.488%
40	14.775	1999	2004	2008	rVB	17512	22493	1.02%	0.109%

41	14.904	2018	2026	2032	rBV4	12840	25665	1.17%	0.124%
42	14.969	2032	2037	2044	rVB2	11886	21661	0.98%	0.105%
43	15.098	2052	2059	2061	rBV2	12643	23724	1.08%	0.115%
44	15.122	2061	2063	2071	rVB5	8894	16213	0.74%	0.079%
45	15.228	2076	2081	2085	rVV3	6450	11243	0.51%	0.054%

46	15.292	2085	2092	2096	rVV3	78007	138526	6.29%	0.671%
47	15.357	2096	2103	2112	rVV	270177	477877	21.70%	2.314%
48	15.445	2112	2118	2121	rVV6	22146	41838	1.90%	0.203%
49	15.481	2121	2124	2126	rVV	24024	33142	1.50%	0.160%
50	15.510	2126	2129	2133	rVV3	46351	62836	2.85%	0.304%

51	15.557	2133	2137	2140	rVV4	9072	16174	0.73%	0.078%
52	15.604	2140	2145	2148	rVV3	19664	35821	1.63%	0.173%
53	15.651	2148	2153	2163	rVV2	66247	120443	5.47%	0.583%
54	15.804	2171	2179	2187	rVV	74597	157735	7.16%	0.764%
55	15.887	2187	2193	2200	rVB3	14225	30815	1.40%	0.149%

56	16.034	2212	2218	2227	rVB5	13935	33836	1.54%	0.164%
57	16.387	2263	2278	2283	rBV2	31085	77535	3.52%	0.375%
58	16.434	2283	2286	2291	rVV2	14216	18629	0.85%	0.090%
59	16.528	2298	2302	2305	rVB	14025	18268	0.83%	0.088%
60	16.569	2305	2309	2312	rBV3	17769	26087	1.18%	0.126%

61	16.604	2312	2315	2319	rVV	15735	21335	0.97%	0.103%
62	16.645	2319	2322	2326	rVB3	13913	20904	0.95%	0.101%
63	16.745	2334	2339	2346	rVB4	20297	38613	1.75%	0.187%
64	16.816	2346	2351	2360	rVV4	21987	62577	2.84%	0.303%
65	16.892	2360	2364	2379	rVB	110943	163424	7.42%	0.791%

66	17.092	2389	2398	2401	rBV	797559	1283611	58.28%	6.216%
67	17.139	2401	2406	2411	rVV	1487637	2202631	100.00%	10.666%
68	17.192	2411	2415	2417	rVV	92152	129707	5.89%	0.628%
69	17.222	2417	2420	2425	rVV	155407	182448	8.28%	0.883%
70	17.275	2425	2429	2436	rVB4	9475	21501	0.98%	0.104%

71	17.516	2462	2470	2474	rBV	50660	89656	4.07%	0.434%
72	17.616	2483	2487	2493	rBV2	19433	28235	1.28%	0.137%
73	17.675	2493	2497	2502	rBV3	18627	28725	1.30%	0.139%
74	17.828	2519	2523	2529	rBV4	11164	19784	0.90%	0.096%
75	17.986	2545	2550	2554	rBV	101359	141178	6.41%	0.684%

76	18.034	2554	2558	2564	rVB	157658	218168	9.90%	1.056%
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 Title : SVOA CALIBRATION

77	18.122	2567	2573	2578	rBV	30539	49001	2.22%	0.237%
78	18.192	2579	2585	2590	rBV2	124820	219571	9.97%	1.063%
79	18.239	2590	2593	2599	rVB2	44404	65590	2.98%	0.318%
80	18.316	2601	2606	2607	rBV3	7784	11869	0.54%	0.057%
81	18.522	2636	2641	2646	rBV	65810	97889	4.44%	0.474%
82	18.569	2646	2649	2654	rVB7	7621	11855	0.54%	0.057%
83	18.663	2663	2665	2668	rVB3	11684	11840	0.54%	0.057%
84	18.757	2678	2681	2686	rVB4	15238	22014	1.00%	0.107%
85	18.951	2710	2714	2718	rBV2	24127	37724	1.71%	0.183%
86	19.086	2733	2737	2744	rBV5	15239	33812	1.54%	0.164%
87	19.216	2752	2759	2765	rBV	825736	1253297	56.90%	6.069%
88	19.286	2769	2771	2775	rBV2	9259	12571	0.57%	0.061%
89	19.328	2775	2778	2781	rVV4	10231	13042	0.59%	0.063%
90	19.469	2798	2802	2807	rBV3	23216	50137	2.28%	0.243%
91	19.563	2812	2818	2819	rVV	217337	311169	14.13%	1.507%
92	19.592	2819	2823	2832	rVV	518674	820308	37.24%	3.972%
93	19.675	2833	2837	2844	rVB2	29047	46206	2.10%	0.224%
94	19.786	2852	2856	2861	rBV	32529	42678	1.94%	0.207%
95	20.116	2908	2912	2916	rVB	69462	99753	4.53%	0.483%
96	20.222	2925	2930	2934	rBV	75232	108773	4.94%	0.527%
97	20.275	2937	2939	2944	rVB2	31667	46018	2.09%	0.223%
98	21.510	3141	3149	3155	rBV2	1105211	2056929	93.39%	9.960%
99	21.557	3155	3157	3168	rVB	136129	221554	10.06%	1.073%
100	24.780	3695	3705	3716	rVB2	640095	1914152	86.90%	9.269%

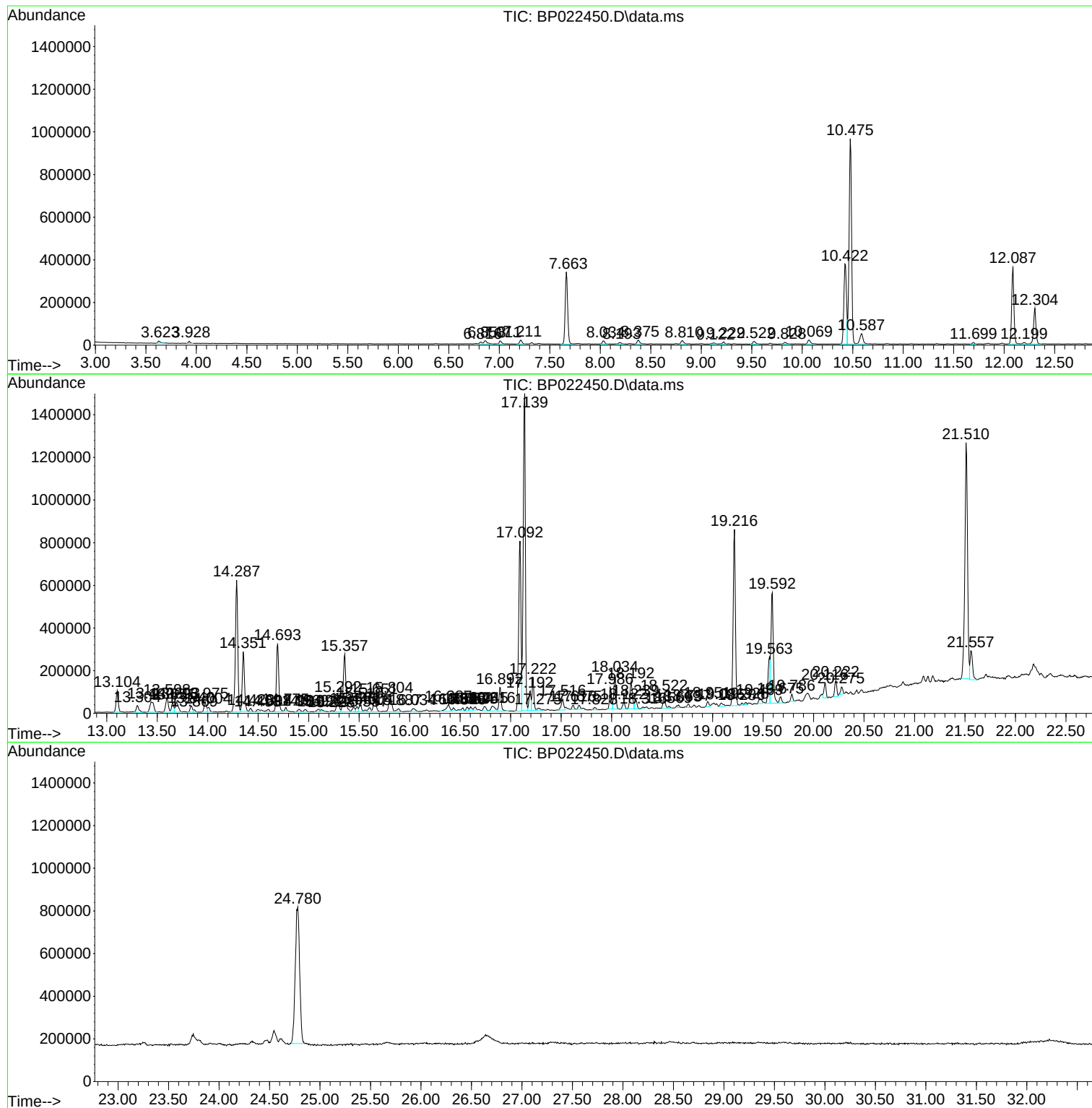
Sum of corrected areas: 20651617

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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Sample : P4264-02MEDL2 20X
Misc :
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Instrument :
BNA_P
ClientSampleId :
DCYN6MEDL2

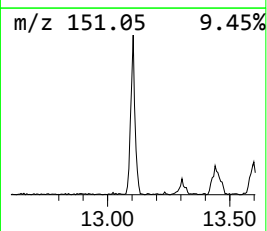
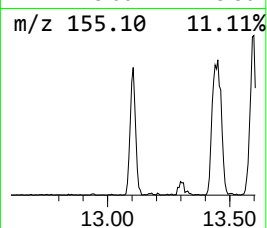
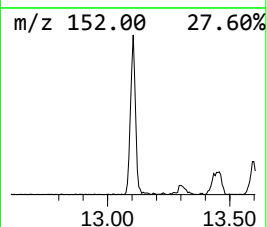
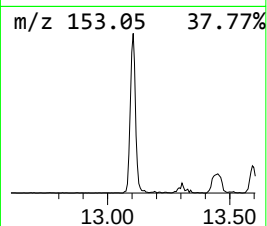
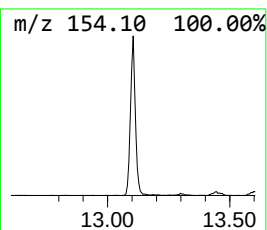
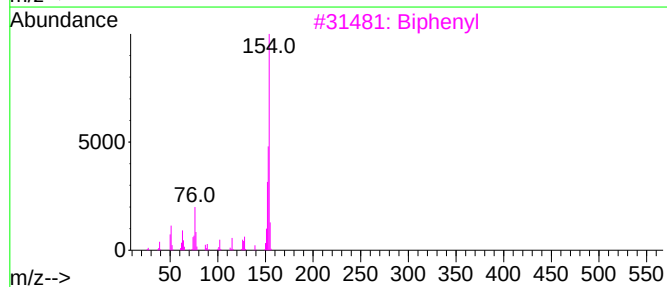
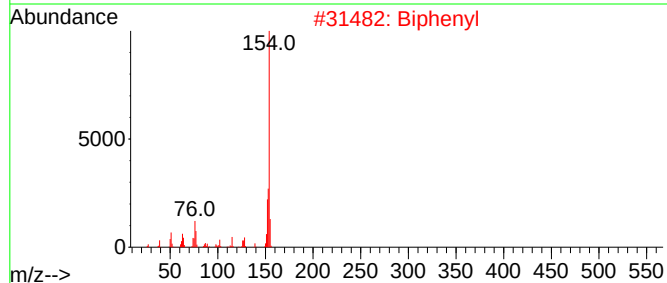
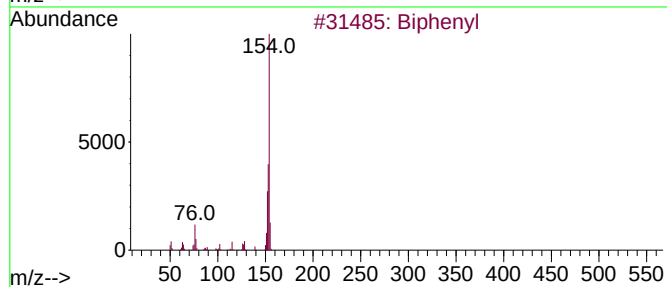
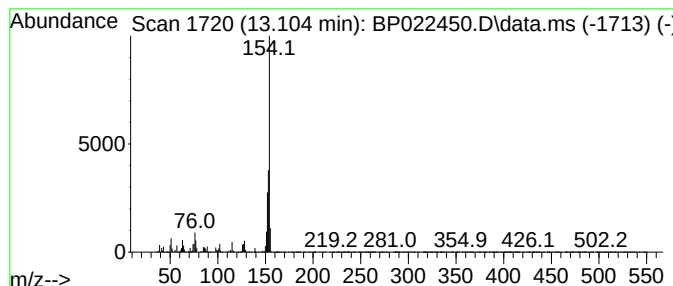
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Biphenyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.104	2.82 ng/ul	146334	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl			154	C12H10	000092-52-4	94
2	Biphenyl			154	C12H10	000092-52-4	93
3	Biphenyl			154	C12H10	000092-52-4	91
4	Naphthalene, 2-ethenyl-			154	C12H10	000827-54-3	91
5	Naphthalene, 2-ethenyl-			154	C12H10	000827-54-3	87



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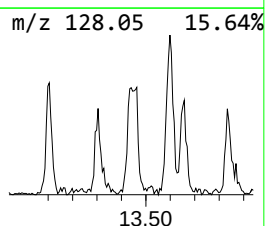
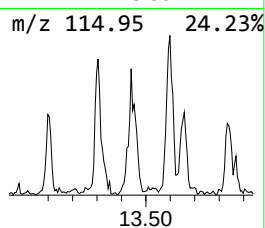
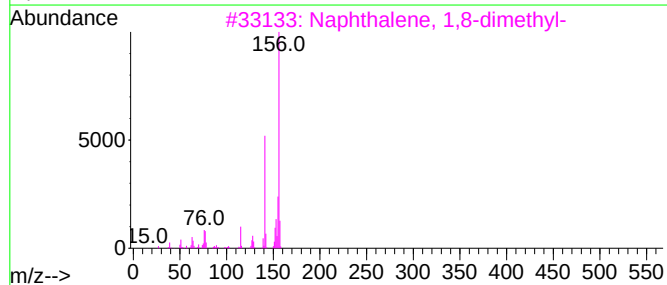
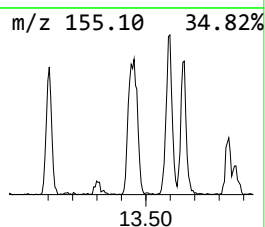
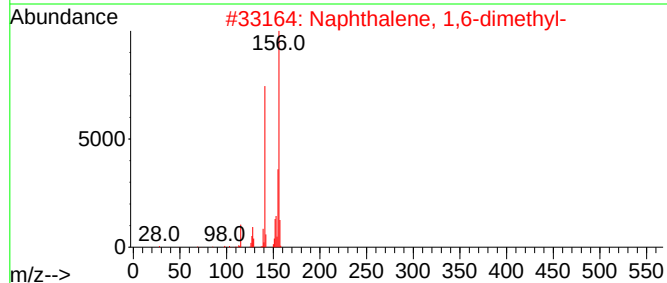
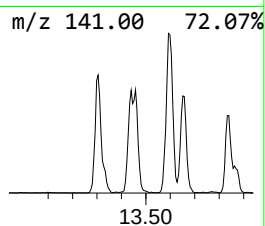
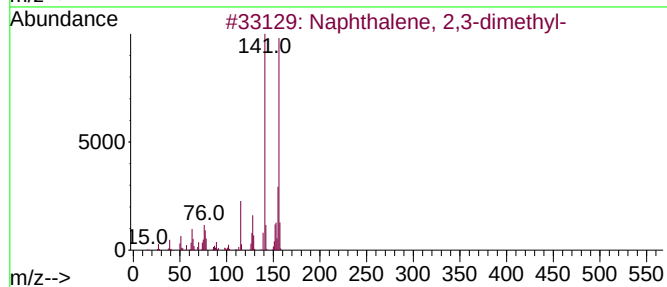
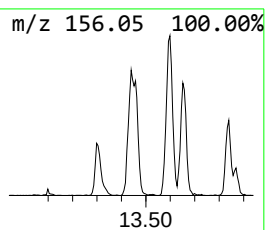
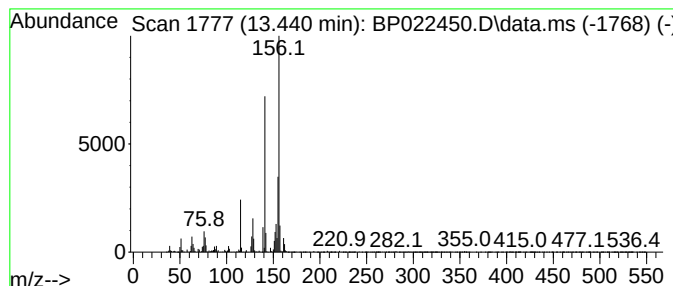
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.440	2.51 ng/ul	130051	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94



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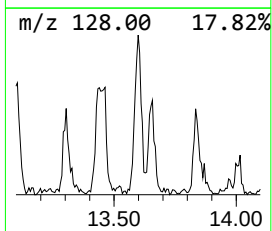
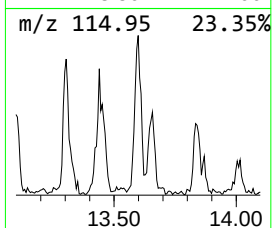
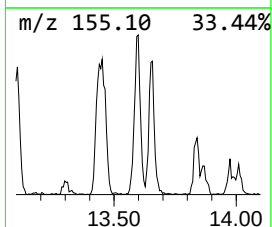
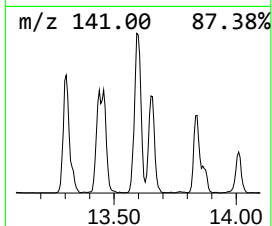
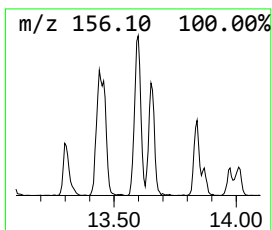
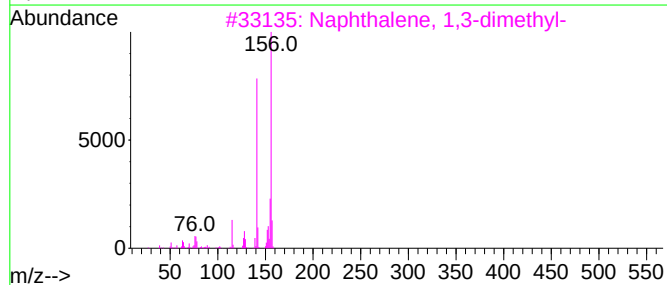
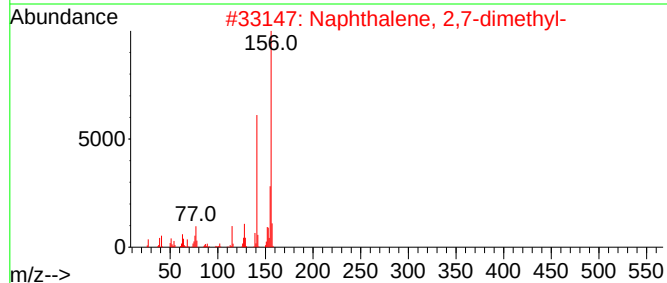
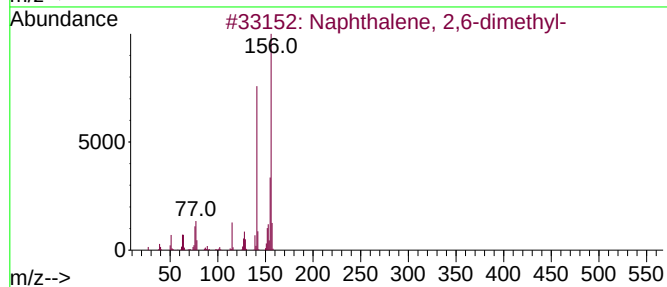
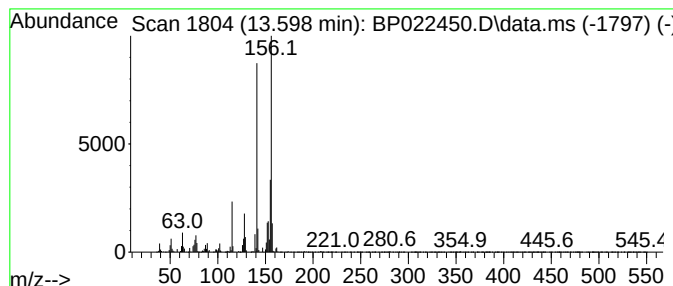
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.599	2.33 ng/ul	120696	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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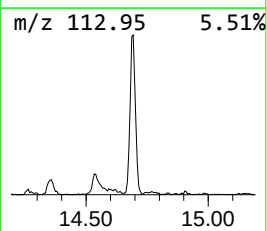
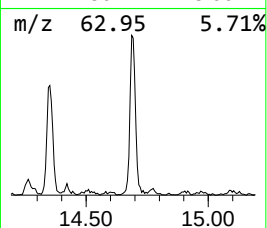
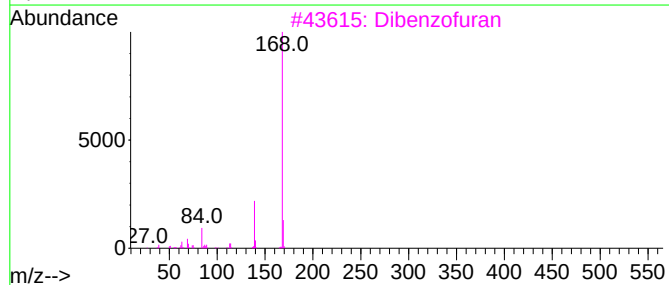
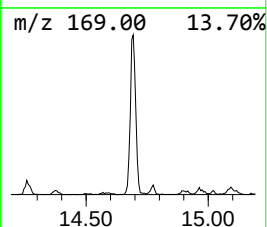
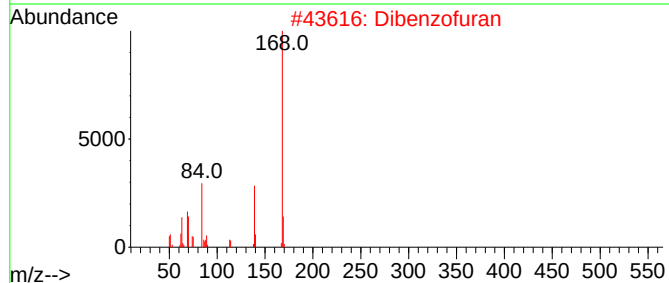
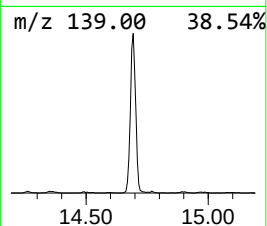
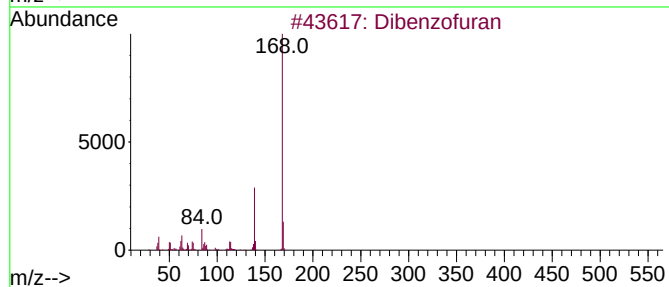
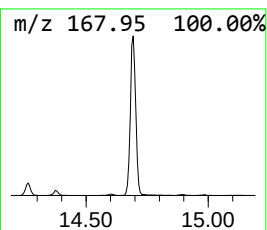
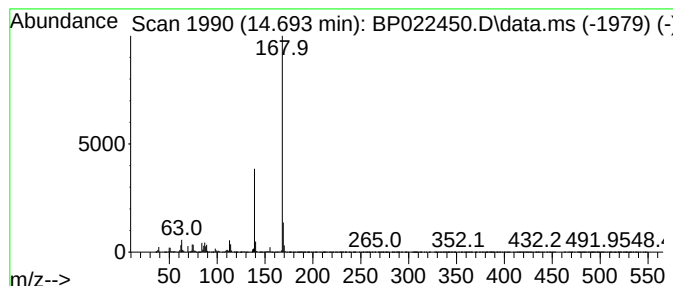
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dibenzo furan Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.693	9.90 ng/ul	513727	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	91
2			Dibenzofuran	168	C12H8O	000132-64-9	78
3			Dibenzofuran	168	C12H8O	000132-64-9	78
4			9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	50
5			1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN6MEDL2

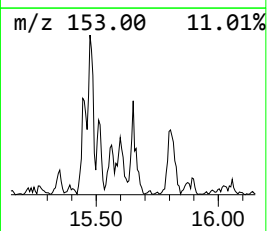
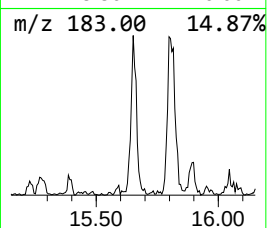
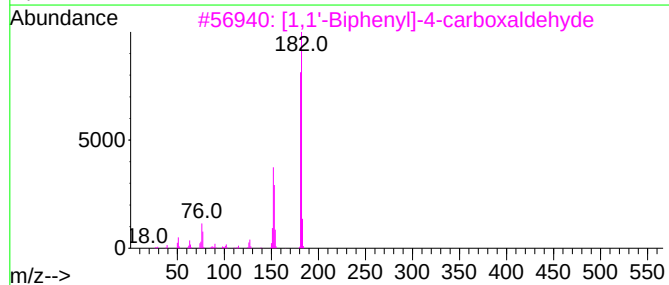
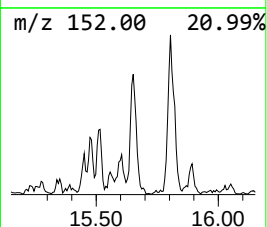
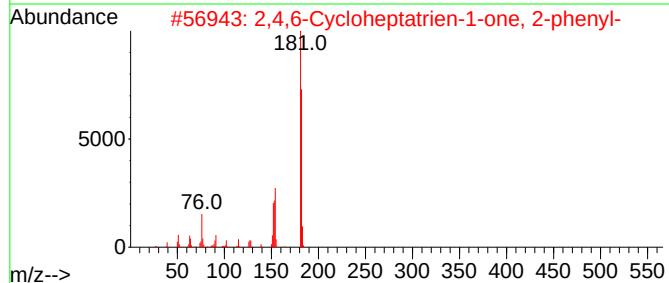
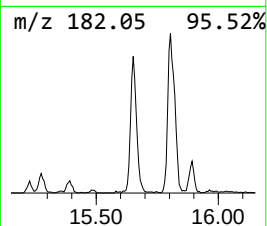
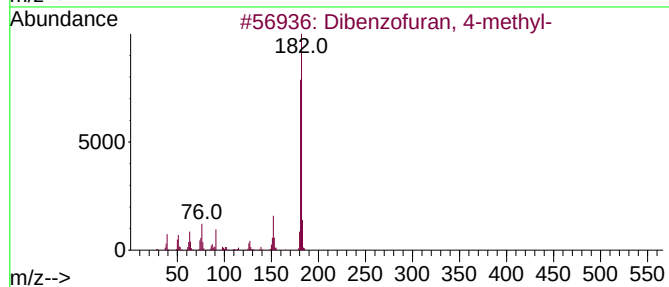
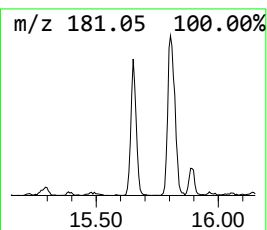
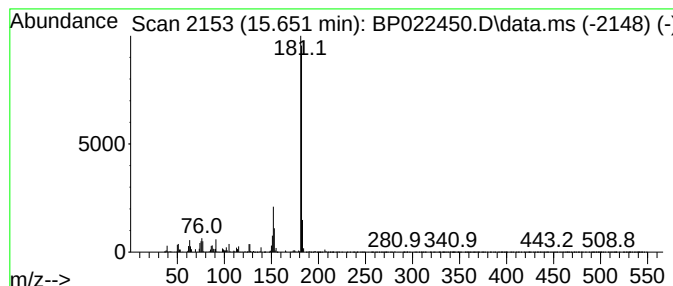
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Dibenzofuran, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	2.32 ng/ul	120443	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
3			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	83
4			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	80
5			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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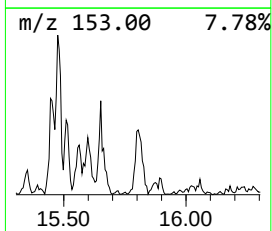
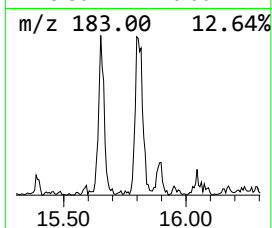
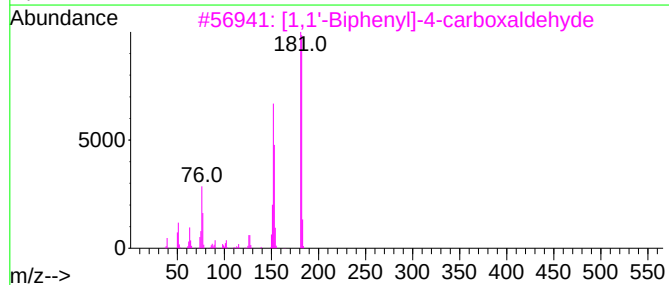
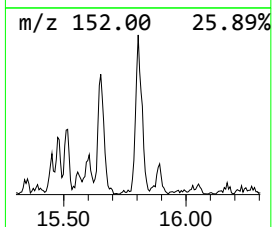
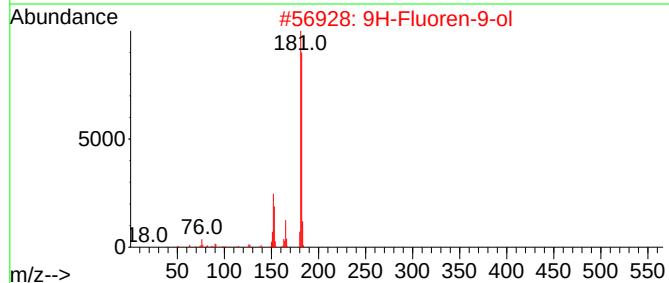
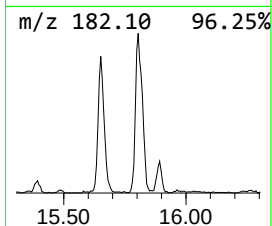
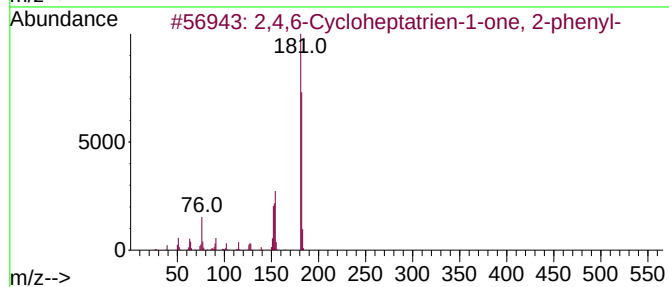
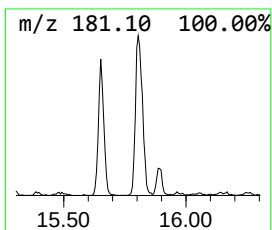
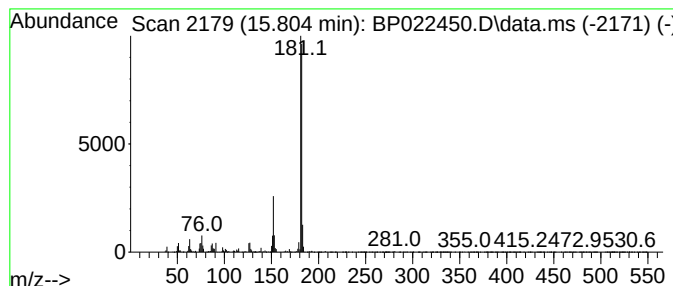
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2,4,6-Cycloheptatrien-1-one... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	2.46 ng/ul	157735	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	91		
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86		
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86		
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86		
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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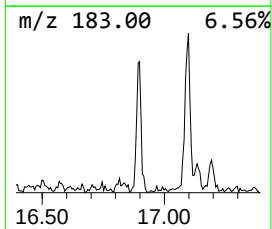
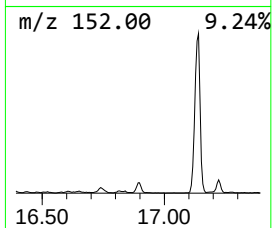
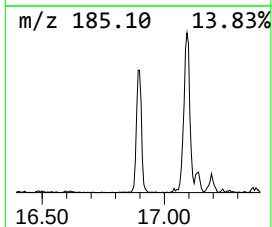
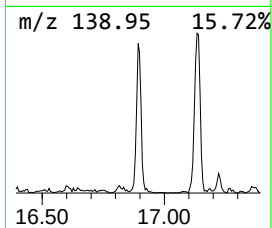
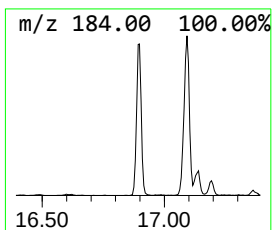
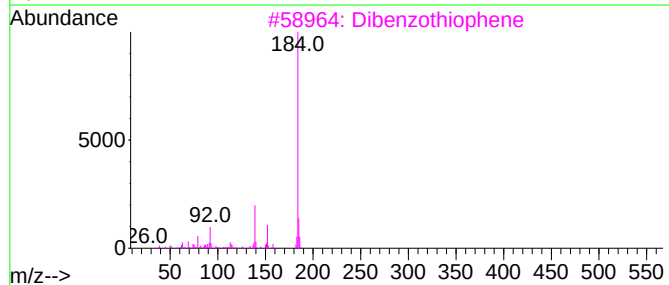
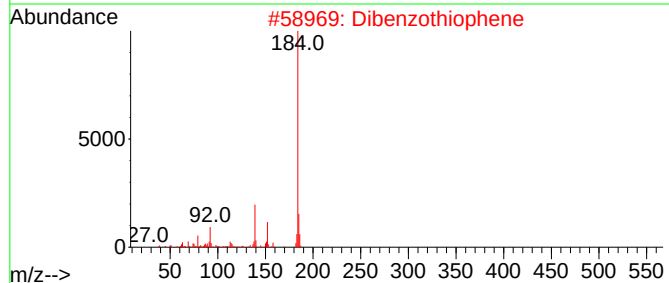
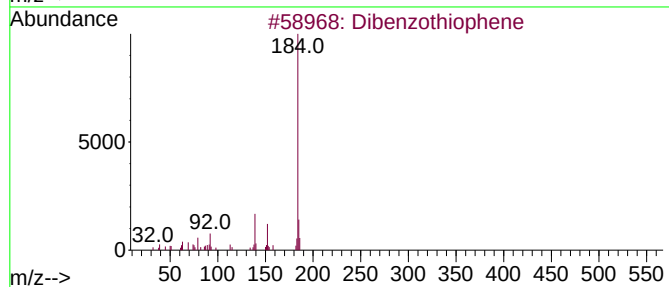
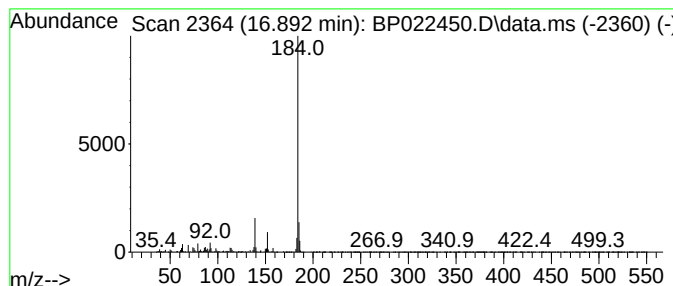
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Dibenzothiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.892	2.55 ng/ul	163424	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	96
2			Dibenzothiophene	184	C12H8S	000132-65-0	95
3			Dibenzothiophene	184	C12H8S	000132-65-0	95
4			Dibenzothiophene	184	C12H8S	000132-65-0	95
5			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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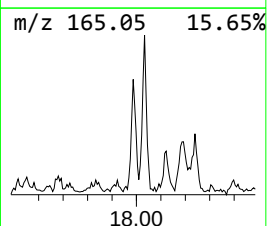
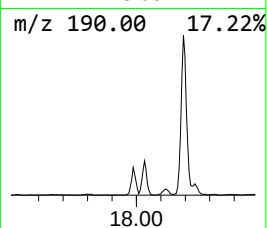
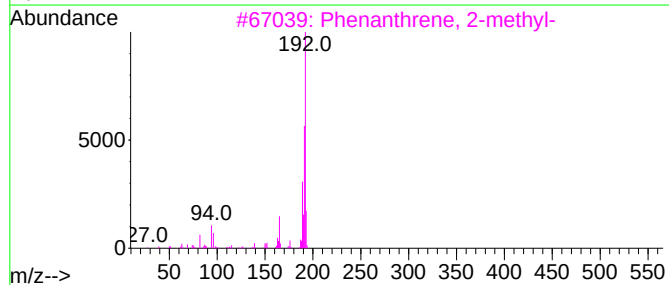
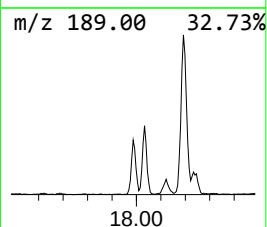
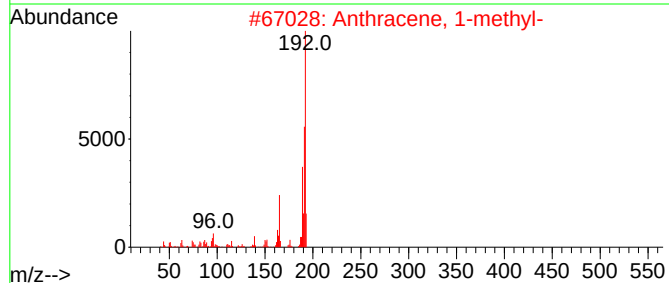
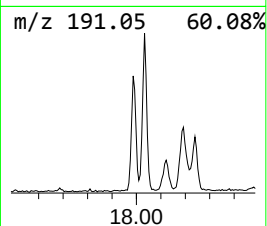
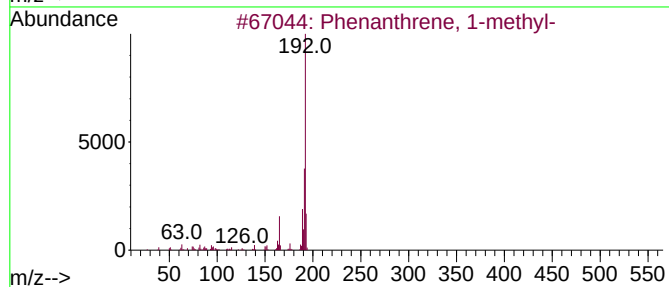
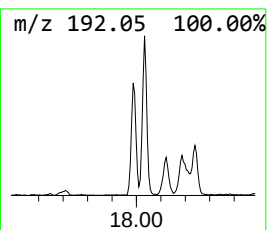
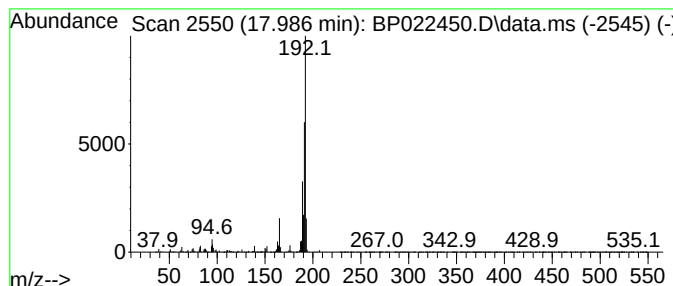
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.986	2.20 ng/ul	141178	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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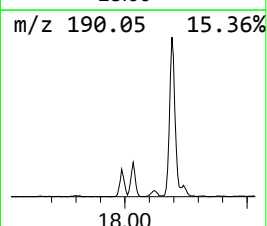
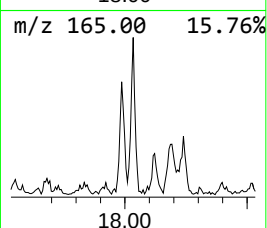
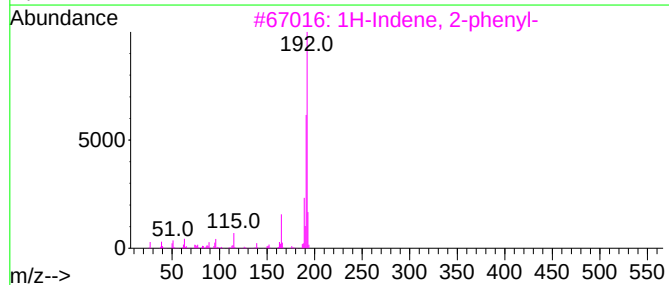
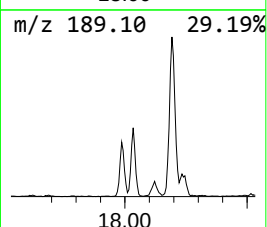
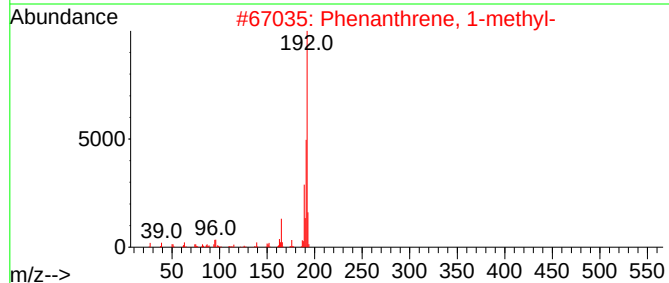
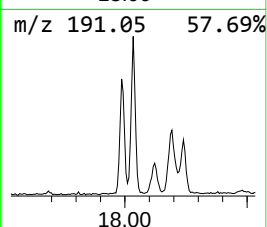
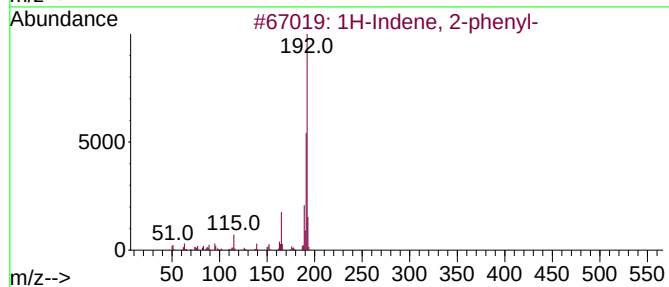
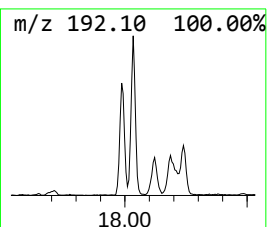
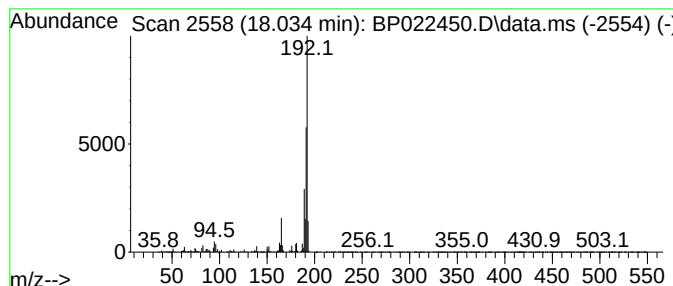
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1H-Indene, 2-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.034	3.40 ng/ul	218168	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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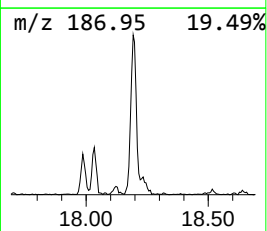
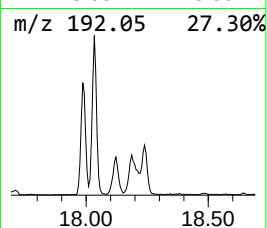
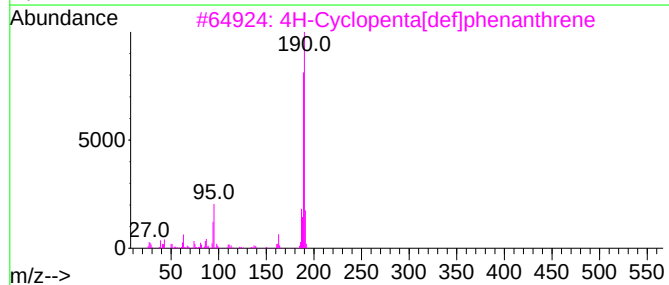
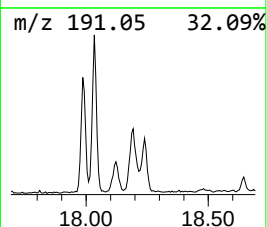
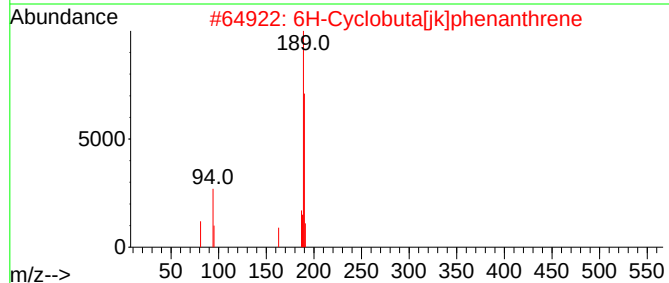
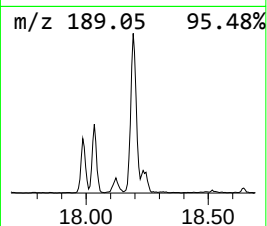
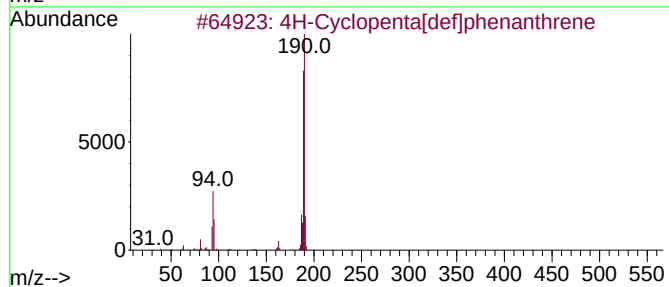
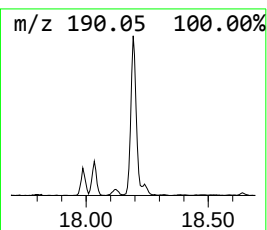
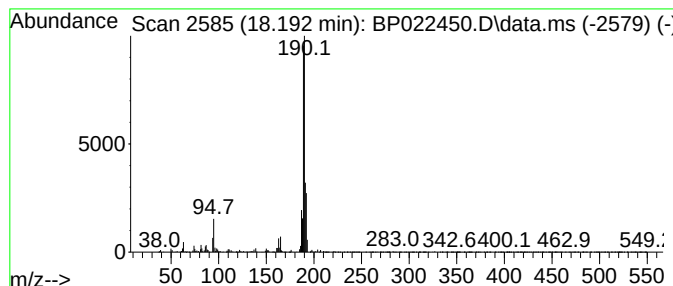
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	3.42 ng/ul	219571	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
4			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	53
5			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
Data File : BP022450.D
Acq On : 17 Oct 2024 18:42
Operator : RC/JU
Sample : P4264-02MEDL2 20X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN6MEDL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Biphenyl	13.104	2.8	ng/ul	146334	3	14.287	1038100	20.0
Naphthalene, 2,...	13.440	2.5	ng/ul	130051	3	14.287	1038100	20.0
Naphthalene, 2,...	13.599	2.3	ng/ul	120696	3	14.287	1038100	20.0
Dibenzo furan	14.693	9.9	ng/ul	513727	3	14.287	1038100	20.0
Dibenzofuran, 4...	15.651	2.3	ng/ul	120443	3	14.287	1038100	20.0
2,4,6-Cyclohept...	15.804	2.5	ng/ul	157735	4	17.092	1283610	20.0
Dibenzothiophene	16.892	2.5	ng/ul	163424	4	17.092	1283610	20.0
Phenanthrene, 1...	17.986	2.2	ng/ul	141178	4	17.092	1283610	20.0
1H-Indene, 2-ph...	18.034	3.4	ng/ul	218168	4	17.092	1283610	20.0
4H-Cyclopenta[d...	18.192	3.4	ng/ul	219571	4	17.092	1283610	20.0