

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
 Data File : BP021171.D
 Acq On : 26 Jul 2024 15:08
 Operator : MA/JU
 Sample : P3347-02DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AW2DL

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021171.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.176	30	32	41	rVB2	13168	19615	0.50%	0.038%
2	3.617	105	107	115	rBV	33238	61018	1.55%	0.118%
3	3.893	151	154	161	rVB	13751	18768	0.48%	0.036%
4	6.746	632	639	642	rBV4	5393	11722	0.30%	0.023%
5	6.875	655	661	675	rBV3	54073	131604	3.34%	0.254%
6	6.993	675	681	691	rVV	159181	254145	6.46%	0.490%
7	7.111	696	701	707	rVB	15059	24076	0.61%	0.046%
8	7.205	709	717	727	rBV	179486	333085	8.46%	0.642%
9	7.287	727	731	738	rVB2	15050	25478	0.65%	0.049%
10	7.640	785	791	798	rVV	761907	1200922	30.52%	2.316%
11	7.705	798	802	820	rVB4	33501	119144	3.03%	0.230%
12	8.005	845	853	859	rVB	38229	66923	1.70%	0.129%
13	8.169	876	881	887	rVB	23188	42428	1.08%	0.082%
14	8.258	891	896	902	rBV4	5878	9845	0.25%	0.019%
15	8.387	912	918	934	rBV	174949	347642	8.83%	0.670%
16	8.722	966	975	982	rBV8	8319	18609	0.47%	0.036%
17	8.805	982	989	1002	rBV	208075	353883	8.99%	0.682%
18	9.105	1029	1040	1044	rBV7	9833	30995	0.79%	0.060%
19	9.299	1067	1073	1078	rVB2	6747	10959	0.28%	0.021%
20	9.516	1101	1110	1121	rBV	162158	297652	7.56%	0.574%
21	9.658	1130	1134	1141	rVV	18132	31429	0.80%	0.061%
22	9.805	1151	1159	1172	rBV4	41735	114617	2.91%	0.221%
23	9.905	1172	1176	1177	rVV3	7460	9414	0.24%	0.018%
24	9.940	1177	1182	1194	rVB6	25155	73742	1.87%	0.142%
25	10.081	1194	1206	1220	rBV	216039	515383	13.10%	0.994%
26	10.210	1223	1228	1236	rVV	21092	47418	1.21%	0.091%
27	10.416	1254	1263	1267	rBV	928404	1732844	44.04%	3.341%
28	10.469	1267	1272	1282	rVB	1664865	2924725	74.33%	5.639%
29	10.581	1284	1291	1299	rBV	210827	390031	9.91%	0.752%
30	10.810	1325	1330	1334	rBV8	10115	22072	0.56%	0.043%
31	10.975	1354	1358	1367	rVB	17387	38087	0.97%	0.073%
32	11.052	1367	1371	1380	rVB4	17797	36159	0.92%	0.070%
33	11.169	1383	1391	1400	rBV7	16039	46244	1.18%	0.089%
34	11.275	1401	1409	1423	rVB2	49340	128456	3.26%	0.248%
35	11.410	1427	1432	1436	rBV8	6162	13736	0.35%	0.026%
36	11.457	1436	1440	1444	rBV3	7281	10885	0.28%	0.021%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Title : SVOA CALIBRATION

37	11.516	1444	1450	1456	rVB3	21651	42398	1.08%	0.082%
38	11.628	1466	1469	1476	rVB9	7897	14261	0.36%	0.027%
39	11.787	1491	1496	1497	rBV5	10896	13485	0.34%	0.026%
40	11.822	1498	1502	1509	rVB2	28719	54229	1.38%	0.105%
41	11.969	1521	1527	1536	rBV3	34281	87639	2.23%	0.169%
42	12.081	1536	1546	1554	rVV	780972	1214009	30.85%	2.341%
43	12.204	1563	1567	1571	rBV2	14335	28251	0.72%	0.054%
44	12.299	1574	1583	1595	rVB	598786	971427	24.69%	1.873%
45	12.434	1603	1606	1609	rBV2	13745	16281	0.41%	0.031%
46	12.475	1610	1613	1618	rVB4	14194	19035	0.48%	0.037%
47	12.622	1634	1638	1641	rVV4	5790	9611	0.24%	0.019%
48	12.740	1648	1658	1659	rBV6	26389	56770	1.44%	0.109%
49	13.110	1713	1721	1727	rVV	99947	199551	5.07%	0.385%
50	13.257	1743	1746	1749	rBV5	11619	10993	0.28%	0.021%
51	13.304	1749	1754	1757	rBV	57145	88379	2.25%	0.170%
52	13.451	1771	1779	1787	rBV3	105072	318957	8.11%	0.615%
53	13.599	1798	1804	1810	rBV	174599	319050	8.11%	0.615%
54	13.657	1810	1814	1817	rVV	118742	184683	4.69%	0.356%
55	13.699	1817	1821	1828	rVB	540056	745968	18.96%	1.438%
56	13.799	1834	1838	1841	rBV3	17550	26156	0.66%	0.050%
57	13.840	1841	1845	1857	rVB2	65528	169025	4.30%	0.326%
58	13.981	1863	1869	1882	rVB	501863	986775	25.08%	1.903%
59	14.222	1905	1910	1914	rBV	157389	213922	5.44%	0.412%
60	14.287	1914	1921	1927	rVV	1465762	2457359	62.45%	4.738%
61	14.357	1927	1933	1941	rVB	2271198	3546010	90.12%	6.837%
62	14.451	1943	1949	1956	rVV	137706	262648	6.67%	0.506%
63	14.581	1967	1971	1974	rBV	53727	85751	2.18%	0.165%
64	14.640	1978	1981	1986	rVV4	40038	72905	1.85%	0.141%
65	14.698	1986	1991	1997	rVB	771320	1174826	29.86%	2.265%
66	14.922	2025	2029	2035	rBV5	23396	38040	0.97%	0.073%
67	15.310	2089	2095	2099	rBV	799439	1103700	28.05%	2.128%
68	15.363	2099	2104	2115	rVB	1397472	1991877	50.62%	3.841%
69	15.487	2117	2125	2129	rBV2	207212	440758	11.20%	0.850%
70	15.581	2138	2141	2145	rBV5	26936	38251	0.97%	0.074%
71	15.622	2145	2148	2151	rVV2	37522	46027	1.17%	0.089%
72	15.669	2152	2156	2161	rVB	85650	118093	3.00%	0.228%
73	15.816	2172	2181	2188	rBV2	106104	263759	6.70%	0.509%
74	16.028	2214	2217	2220	rBV2	22722	28801	0.73%	0.056%
75	16.075	2220	2225	2234	rVB	196193	336968	8.56%	0.650%
76	16.298	2260	2263	2267	rVB2	34177	47378	1.20%	0.091%

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Integration Parameters: LSCINT.P

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

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 Peak separation: 5

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77	16.387	2270	2278	2286	rBV7	124073	350514	8.91%	0.676%
78	16.769	2338	2343	2352	rBV	230789	343185	8.72%	0.662%
79	16.916	2363	2368	2374	rBV	171000	269198	6.84%	0.519%
80	17.022	2383	2386	2393	rVB7	24926	43695	1.11%	0.084%
81	17.110	2394	2401	2405	rBV	1838478	2898496	73.66%	5.589%
82	17.157	2405	2409	2414	rVV	1870323	2645688	67.24%	5.101%
83	17.210	2414	2418	2421	rVV2	844967	1258112	31.97%	2.426%
84	17.245	2421	2424	2429	rVB	615303	852105	21.65%	1.643%
85	17.551	2466	2476	2482	rBV	2435412	3934923	100.00%	7.587%
86	17.722	2502	2505	2513	rBV3	63943	115458	2.93%	0.223%
87	18.034	2554	2558	2562	rBV3	142247	230535	5.86%	0.445%
88	18.069	2562	2564	2569	rVB2	111060	151164	3.84%	0.291%
89	18.222	2585	2590	2604	rVB2	163963	342949	8.72%	0.661%
90	18.345	2607	2611	2615	rVV	71528	116789	2.97%	0.225%
91	18.392	2616	2619	2622	rVB	71150	88780	2.26%	0.171%
92	18.604	2653	2655	2664	rVB	68698	104674	2.66%	0.202%
93	19.039	2726	2729	2734	rBV	53643	71098	1.81%	0.137%
94	19.245	2760	2764	2771	rVB	342899	479534	12.19%	0.925%
95	19.375	2782	2786	2792	rBV3	95718	176335	4.48%	0.340%
96	19.592	2818	2823	2832	rBV2	1097869	1724138	43.82%	3.324%
97	20.122	2910	2913	2921	rBV	105799	181441	4.61%	0.350%
98	21.557	3152	3157	3169	rVB2	1856835	3190644	81.09%	6.152%
99	24.651	3677	3683	3693	rVB	619852	1411480	35.87%	2.722%
100	24.886	3714	3723	3739	rVB2	1207999	3523226	89.54%	6.793%

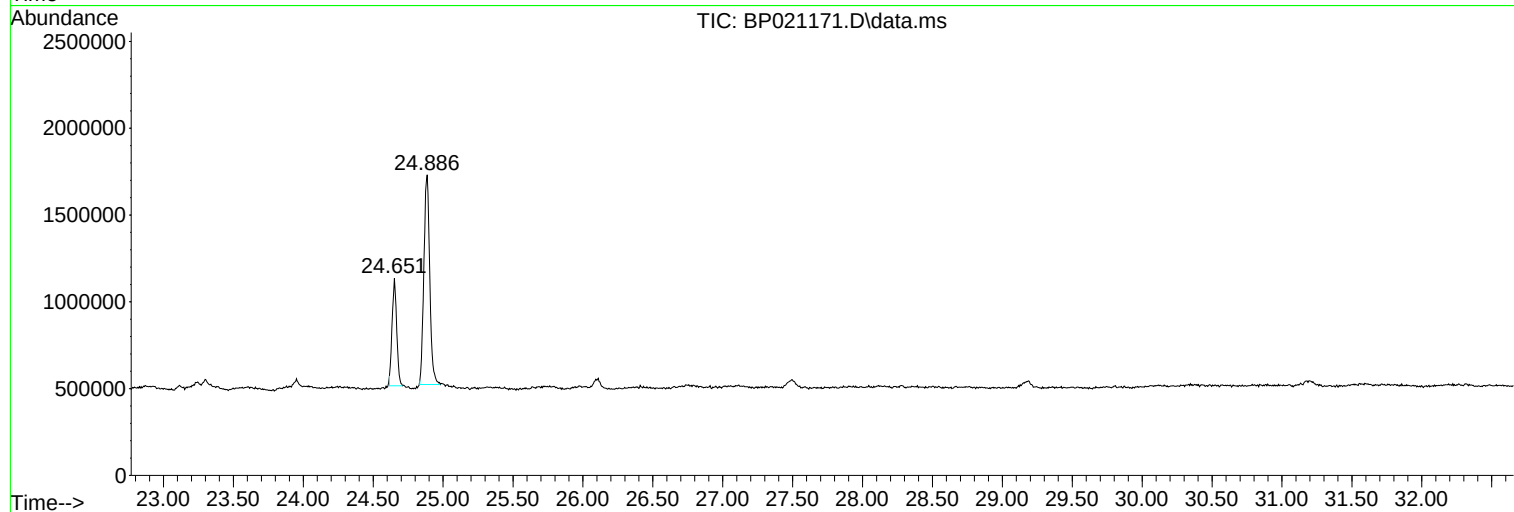
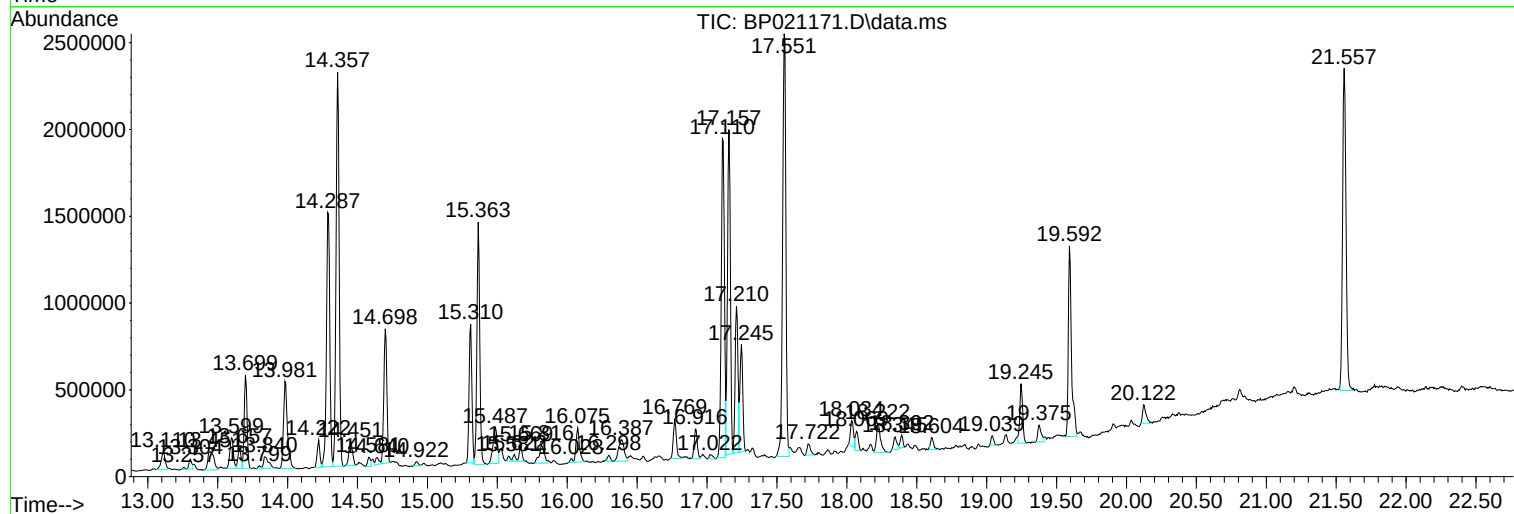
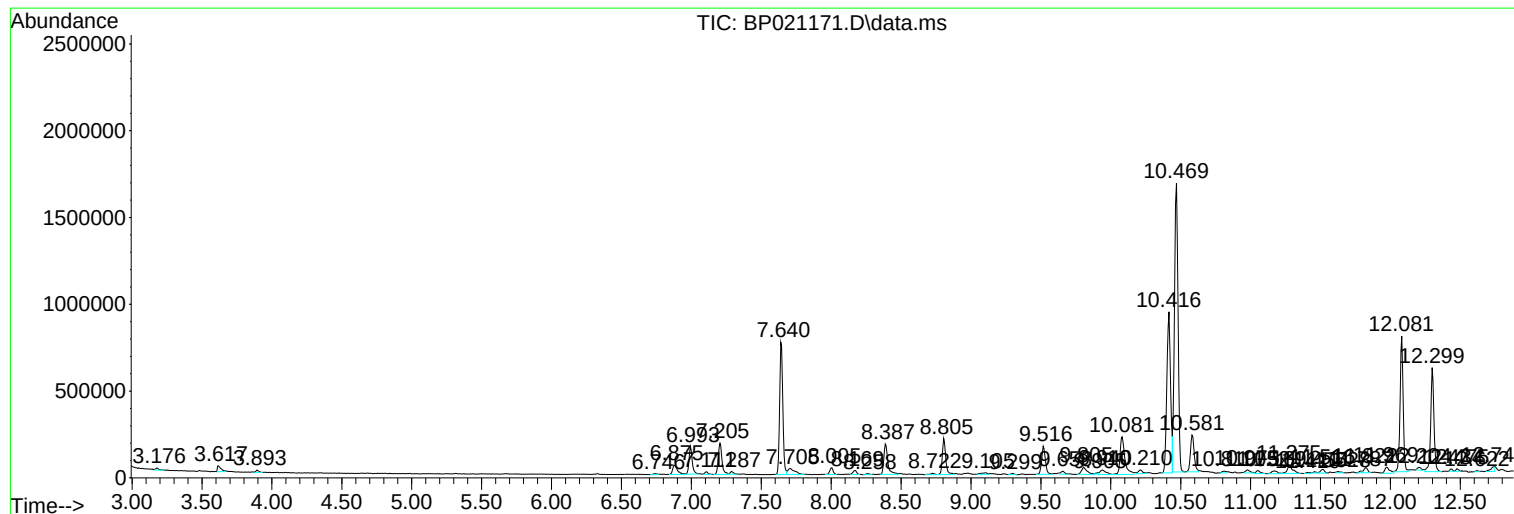
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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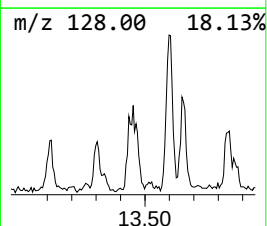
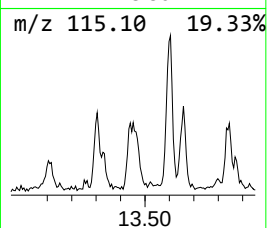
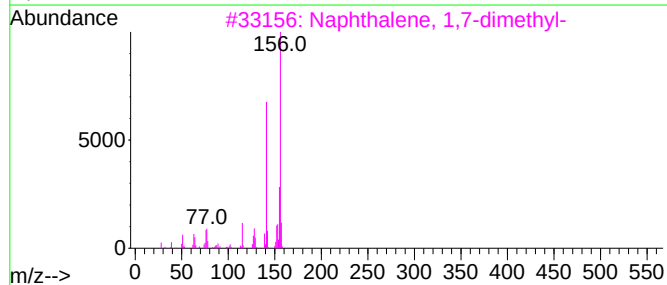
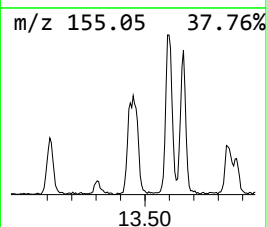
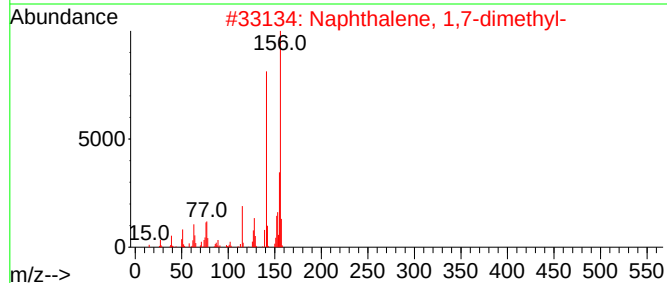
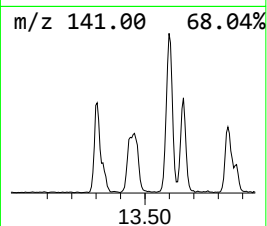
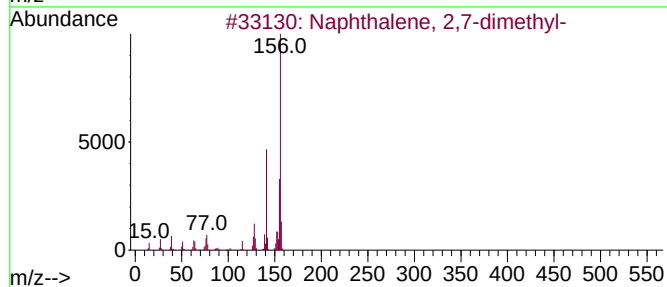
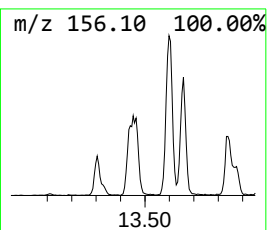
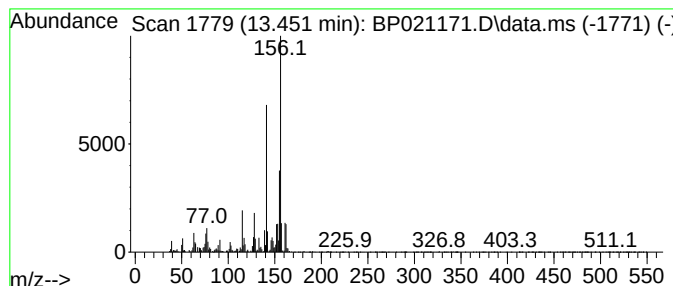
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Naphthalene, 2,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.451	2.60 ng/ul	318957	Acenaphthene-d10	14.293

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96



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BNA_P
ClientSampleId :
C0AW2DL

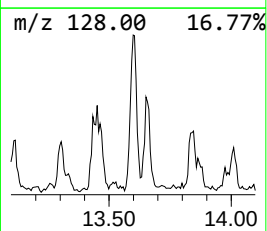
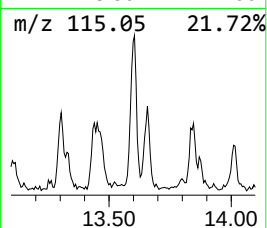
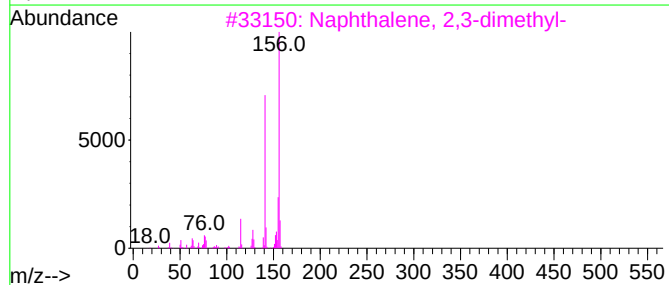
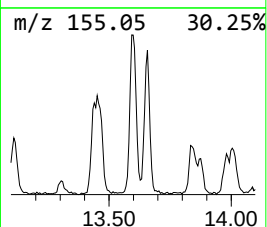
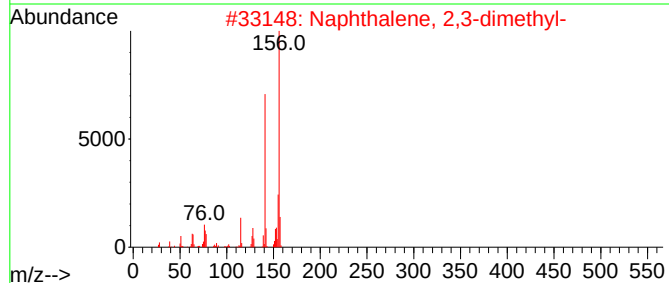
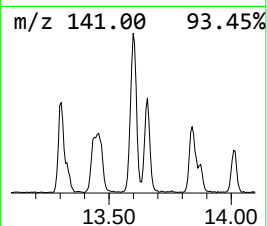
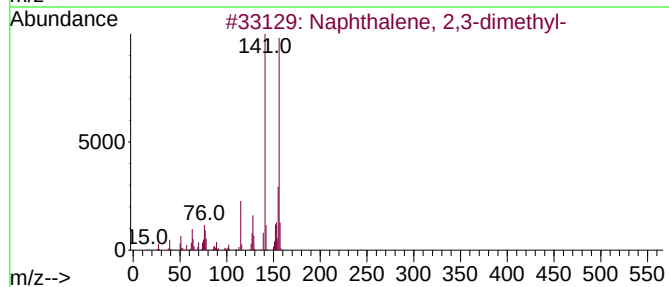
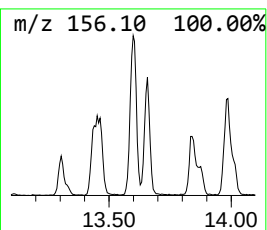
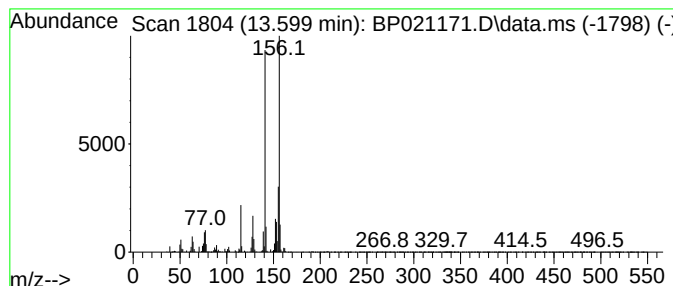
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.599	2.60 ng/ul	319050	Acenaphthene-d10	14.293

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



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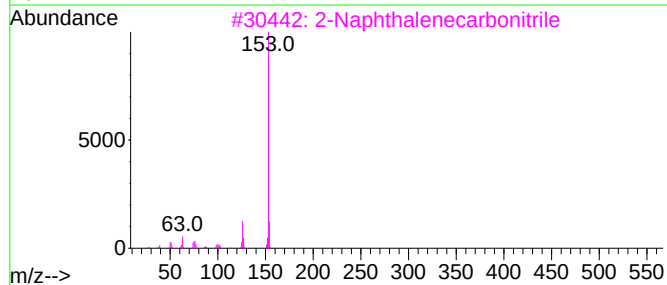
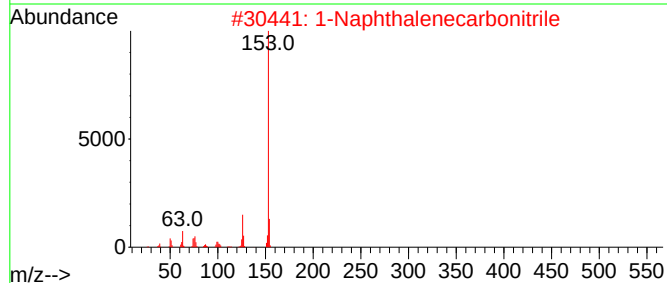
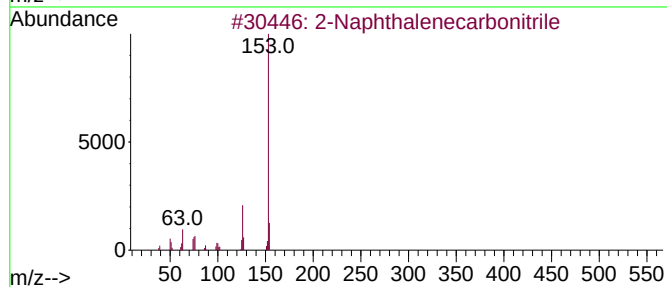
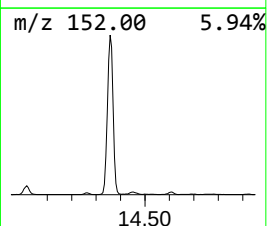
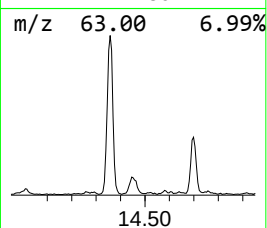
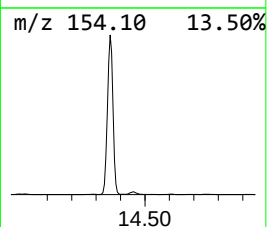
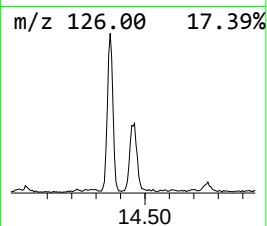
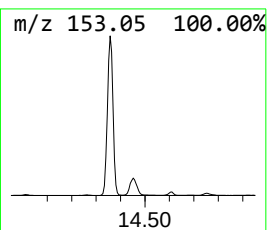
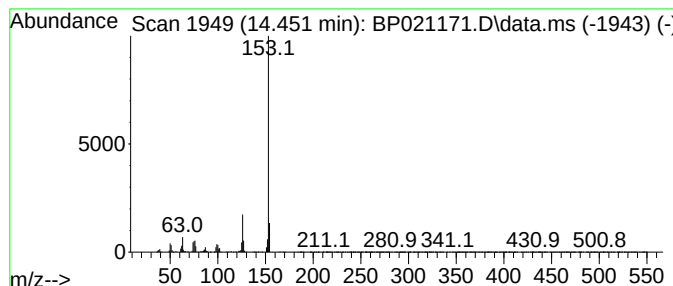
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Naphthalenecarbonitrile Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.451	2.14 ng/ul	262648	Acenaphthene-d10	14.293

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97		
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	97		
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	95		
4	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	94		
5	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021171.D
Acq On : 26 Jul 2024 15:08
Operator : MA/JU
Sample : P3347-02DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW2DL

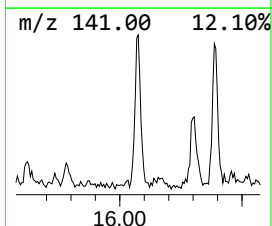
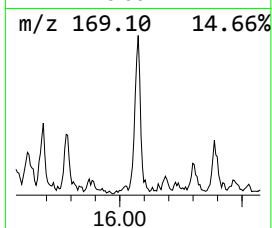
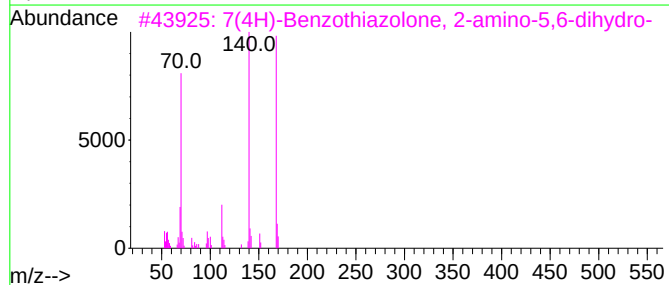
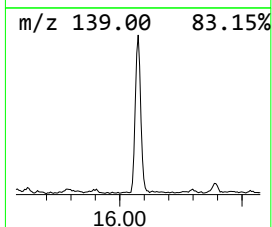
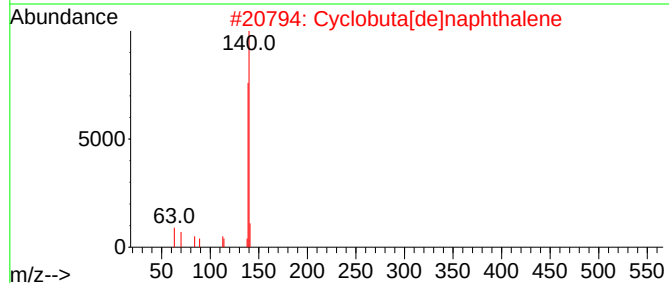
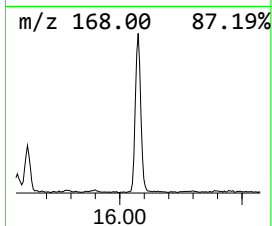
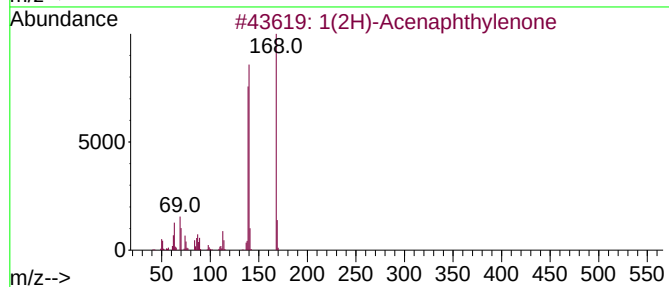
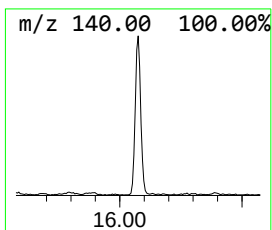
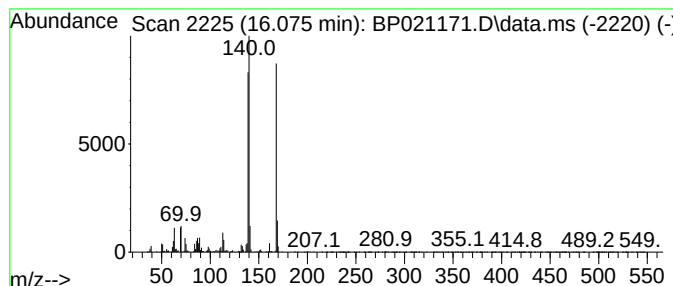
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1(2H)-Acenaphthylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	2.33 ng/ul	336968	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1(2H)-Acenaphthylene	168	C12H8O	002235-15-6	96
2			Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	60
3			7(4H)-Benzothiazolone, 2-amino-5...	168	C7H8N2OS	017583-10-7	50
4			Alanylalanylalanine, N,N',N"-tri...	411	C20H33N3O6	1000329-34-4	50
5			Ethyl maltol	140	C7H8O3	004940-11-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021171.D
Acq On : 26 Jul 2024 15:08
Operator : MA/JU
Sample : P3347-02DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW2DL

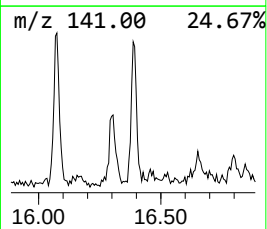
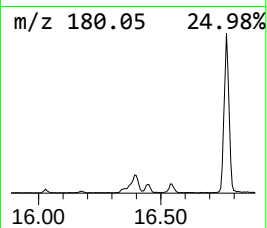
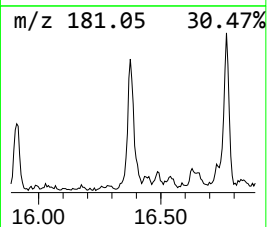
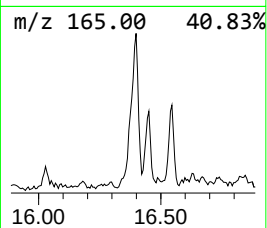
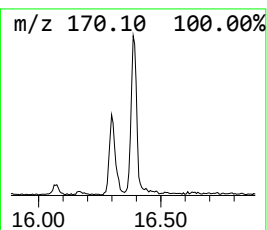
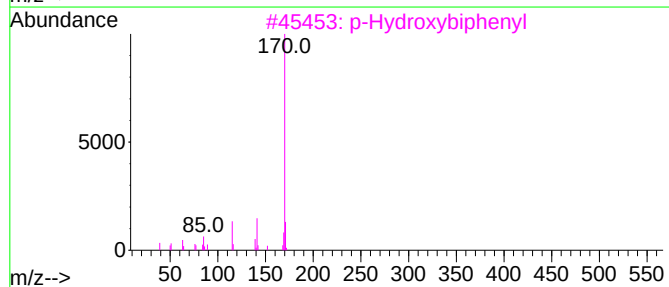
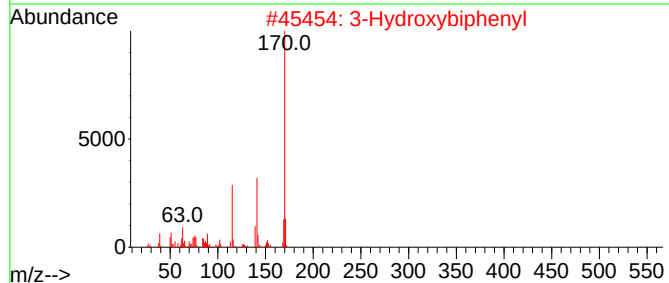
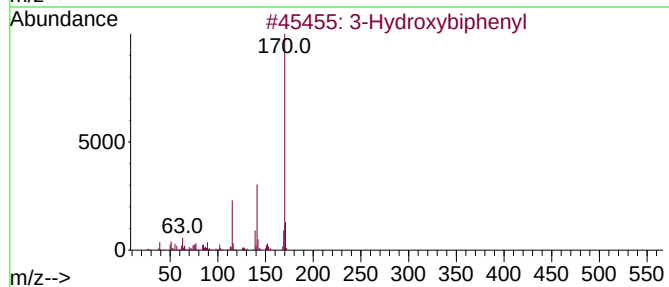
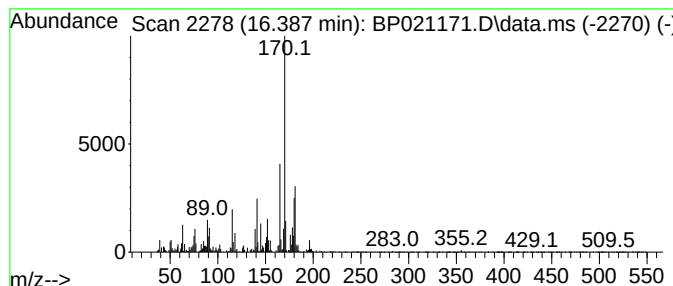
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 3-Hydroxybiphenyl Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.387	2.42 ng/ul	350514	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Hydroxybiphenyl	170	C12H10O	000580-51-8	55
2			3-Hydroxybiphenyl	170	C12H10O	000580-51-8	50
3			p-Hydroxybiphenyl	170	C12H10O	000092-69-3	49
4			p-Hydroxybiphenyl	170	C12H10O	000092-69-3	49
5			3-Hydroxybiphenyl	170	C12H10O	000580-51-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021171.D
Acq On : 26 Jul 2024 15:08
Operator : MA/JU
Sample : P3347-02DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW2DL

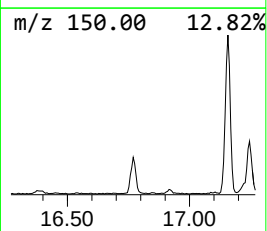
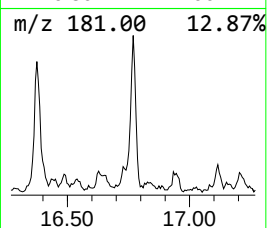
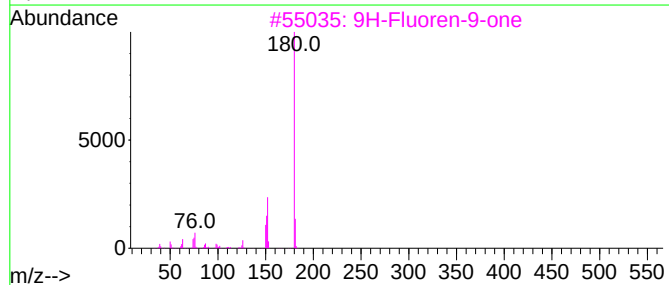
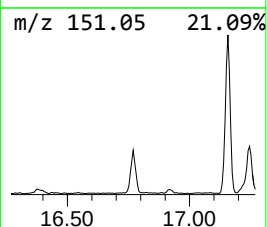
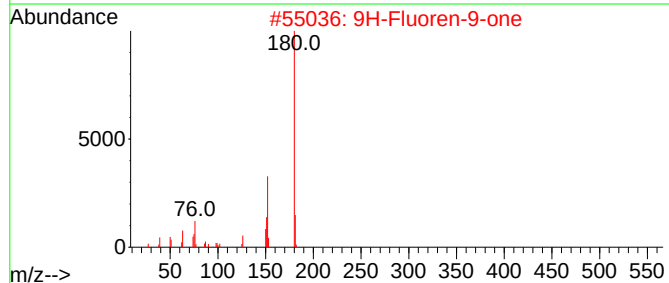
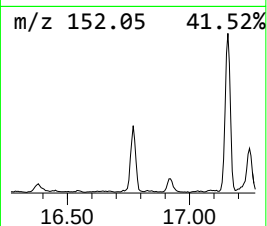
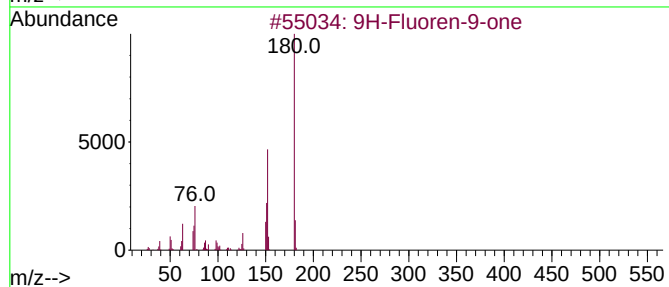
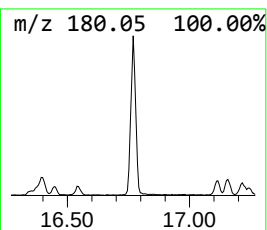
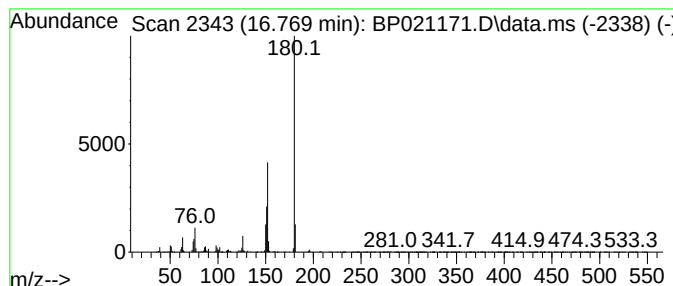
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9H-Fluoren-9-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.769	2.37 ng/ul	343185	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one			180	C13H8O	000486-25-9	95
2	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
3	9H-Fluoren-9-one			180	C13H8O	000486-25-9	94
4	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91
5	Benzo[h]cinnoline			180	C12H8N2	000230-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021171.D
Acq On : 26 Jul 2024 15:08
Operator : MA/JU
Sample : P3347-02DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW2DL

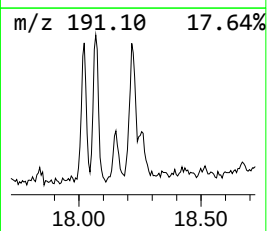
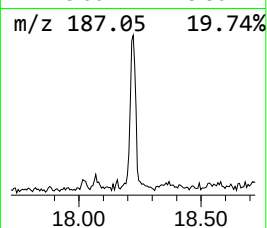
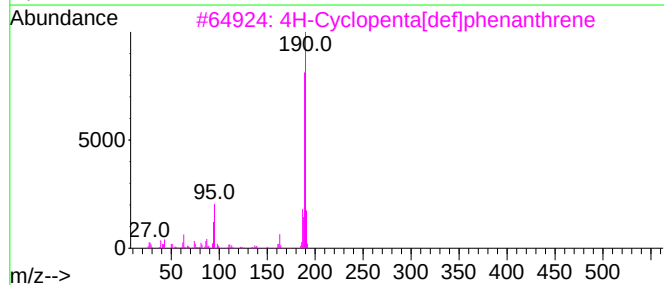
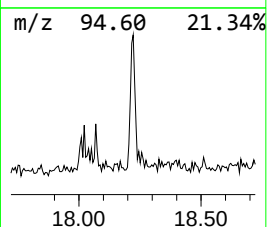
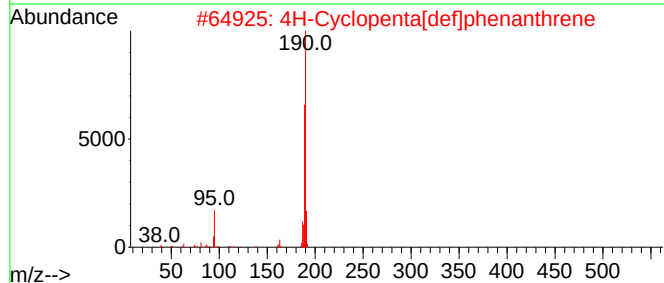
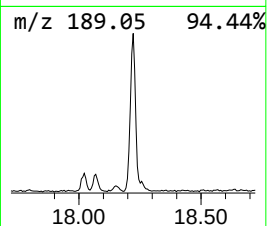
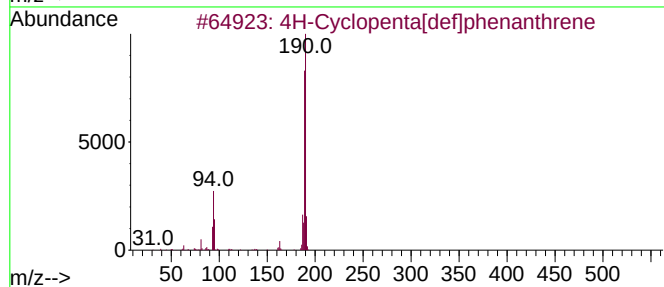
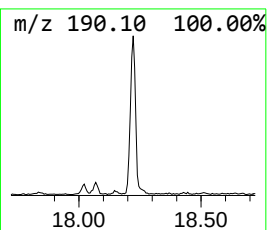
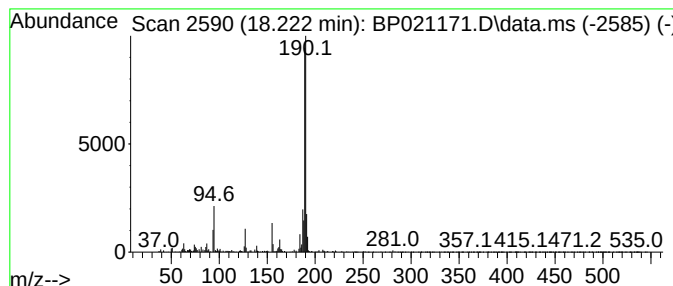
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	2.37 ng/ul	342949	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	80
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	64
5			Phenol, 4-(2-thienylmethyl)-	190	C11H10O5	091680-55-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021171.D
Acq On : 26 Jul 2024 15:08
Operator : MA/JU
Sample : P3347-02DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
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Naphthalene, 2,...	13.599	2.6	ng/ul	319050	3	14.293	2457360	20.0
2-Naphthaleneca...	14.451	2.1	ng/ul	262648	3	14.293	2457360	20.0
1(2H)-Acenaphth...	16.075	2.3	ng/ul	336968	4	17.110	2898500	20.0
3-Hydroxybiphenyl	16.387	2.4	ng/ul	350514	4	17.110	2898500	20.0
9H-Fluoren-9-one	16.769	2.4	ng/ul	343185	4	17.110	2898500	20.0
4H-Cyclopenta[d...	18.222	2.4	ng/ul	342949	4	17.110	2898500	20.0