LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BHBX9

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_N\Methods\SFAM-EPA-BN052924.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BN031828.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.452	72	79	89	rVB	112024	196515	2.45%	0.119%
2	3.593	100	103	115	rBV2	454431	664457	8.28%	0.402%
3	3.693	116	120	128	rVB	330489	437123	5.45%	0.264%
4	4.893	319	324	331	rVV2	137639	234561	2.92%	0.142%
5	5.675	452	457	465	rVV	445734	779512	9.72%	0.471%
6	6.658	615	624	630	rBV3	85524	192695	2.40%	0.116%
7	6.846	646	656	662	rVV	1665593	2910198	36.28%	1.759%
8	7.011	670	684	696	rBV	1353209	2239148	27.91%	1.353%
9	7.205	710	717	724	rVB	1701485	2695913	33.61%	1.629%
10	7.669	787	796	803	rVV	1620257	2622652	32.69%	1.585%
11	8.375	910	916	924	rBV	1787660	2916967	36.36%	1.763%
12	8.828	985	993	999	rVV	1556988	2595087	32.35%	1.568%
13	9.393	1083	1089	1099	rVB	206389	350186	4.37%	0.212%
14	9.546	1106	1115	1126	rBV	1250416	2201348	27.44%	1.330%
15	10.075	1193	1205	1215	rVB	2020219	3470617	43.26%	2.097%
16	10.452	1260	1269	1275	rBV	2137169	3944437	49.17%	2.383%
17	10.505	1275	1278	1284	rVV	138958	220803	2.75%	0.133%
18	10.599	1284	1294	1314	rVV	556593	1246063	15.53%	0.753%
19	11.199	1390	1396	1407	rVV	321005	637721	7.95%	0.385%
20	11.881	1506	1512	1518	rVV	125610	204126	2.54%	0.123%
21	12.122	1547	1553	1561	rVB	199627	352895	4.40%	0.213%
22	12.340	1584	1590	1597	rBV	132856	238626	2.97%	0.144%
23	13.140	1720	1726	1733	rVB2	224624	345812	4.31%	0.209%
24	13.634	1805	1810	1815	rVV	239514	393925	4.91%	0.238%
25	13.687	1815	1819	1821	rBV	151398	190704	2.38%	0.115%
26	13.728	1821	1826	1832	rVB	3235353	4657433	58.06%	2.814%
27	14.004	1867	1873	1879	rBV2	3976500	6265564	78.10%	3.786%
28	14.216	1905	1909	1915	rVB	186017	248223	3.09%	0.150%
29	14.316	1916	1926	1933	rBV	3148152	5312903	66.23%	3.210%
30	14.528	1955	1962	1970	rBV	1214761	2053344	25.60%	1.241%
31	14.681	1982	1988	1990	rVV	222173	334810	4.17%	0.202%
32	14.716	1990	1994	2001	rVV	285333	495596	6.18%	0.299%
33	14.793	2002	2007	2011	rVB	185109	260332	3.25%	0.157%
34	14.934	2024	2031	2035	rBV4	119360	232526	2.90%	0.141%
35	14.987	2036	2040	2044	rVB	190558	252986	3.15%	0.153%
36	15.175	2068	2072	2078	rVB2	133983	202859	2.53%	0.123%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

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_N\Methods\SFAM-EPA-BN052924.MA.M
Title : SVOA CALIBRATION

37	15.310	2088	2095	2101	rVV	5016475	7419044	92.48%	4.483%
38	15.363	2101	2104	2107	rVV3	154531	224474	2.80%	0.136%
39	15.404	2107	2111	2112	rVV	172278	222397	2.77%	0.134%
40	15.440	2112	2117	2122	rVV	1036281	1467264	18.29%	0.887%
41	15.657	2149	2154	2160	rVV2	147776	255847	3.19%	0.155%
42	15.804	2172	2179	2187	rVB4	77273	219953	2.74%	0.133%
43	16.040	2211	2219	2226	rVB5	271984	657487	8.20%	0.397%
44	16.375	2267	2276	2279	rBV4	127639	324615	4.05%	0.196%
45	16.416	2279	2283	2289	rVB	320045	456146	5.69%	0.276%
46	16.669	2322	2326	2331	rVV	749987	1026336	12.79%	0.620%
47	16.722	2331	2335	2342	rVB2	367417	593564	7.40%	0.359%
48	16.828	2350	2353	2358	rVV2	132113	247084	3.08%	0.149%
49	16.881	2358	2362	2371	rVB2	243241	398053	4.96%	0.241%
50	17.063	2384	2393	2397	rBV	3243956	4772127	59.49%	2.884%
51	17.104	2397	2400	2405	rVV	2677852	3694948	46.06%	2.233%
52	17.163	2405	2410	2414	rVV2	4513336	6592025	82.17%	3.983%
53	17.198	2414	2416	2420	rVV	469793	527232	6.57%	0.319%
54	17.257	2420	2426	2430	rVV2	208956	416941	5.20%	0.252%
55	17.345	2436	2441	2443	rVV2	167048	325239	4.05%	0.197%
56	17.375	2443	2446	2451	rVB3	297801	437634	5.46%	0.264%
57	17.469	2458	2462	2471	rVB	194980	263261	3.28%	0.159%
58	17.645	2488	2492	2500	rVB2	361445	554563	6.91%	0.335%
59	17.939	2537	2542	2544	rVV	821085	1218193	15.19%	0.736%
60	17.975	2544	2548	2556	rVB2	1548678	3579615	44.62%	2.163%
61	18.128	2569	2574	2578	rBV3	790659	1343640	16.75%	0.812%
62	18.169	2578	2581	2589	rVB	480142	667571	8.32%	0.403%
63	18.257	2592	2596	2599	rBV4	259877	387434	4.83%	0.234%
64	18.334	2606	2609	2614	rVB2	216800	305934	3.81%	0.185%
65	18.445	2623	2628	2631	rBV	567459	809068	10.09%	0.489%
66	18.486	2631	2635	2645	rVB2	499654	838037	10.45%	0.506%
67	18.686	2666	2669	2674	rVV	291907	402373	5.02%	0.243%
68	18.739	2674	2678	2682	rVV	416931	568619	7.09%	0.344%
69	18.781	2682	2685	2689	rVB	297540	369596	4.61%	0.223%
70	18.863	2694	2699	2703	rBV	689526	1058333	13.19%	0.640%
71	18.910	2703	2707	2712	rVV4	261940	554161	6.91%	0.335%
72	18.951	2712	2714	2718	rVB	218791	255284	3.18%	0.154%
73	19.004	2718	2723	2729	rBV2	398008	782917	9.76%	0.473%
74	19.128	2738	2744	2751	rVB	5160726	7065142	88.07%	4.269%
75	19.192	2751	2755	2758	rBV	319067	381897	4.76%	0.231%
76	19.251	2762	2765	2773	rVB4	433661	779472	9.72%	0.471%

LSC Area Percent Report

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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

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

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Title : SVOA CALIBRATION

77	19.463	2795	2801	2803	rBV3	4742364	6845541	85.33%	4.137%
78	19.492	2803	2806	2814	rVB	4188367	5810073	72.43%	3.511%
79	19.563	2814	2818	2824	rVB2	367473	569513	7.10%	0.344%
80	19.816	2856	2861	2871	rBV3	449605	1108264	13.82%	0.670%
81	19.951	2879	2884	2886	rBV4	338597	577099	7.19%	0.349%
82	19.986	2887	2890	2893	rVB	660275	827588	10.32%	0.500%
83	20.086	2904	2907	2911	rBV	422650	537975	6.71%	0.325%
84	20.145	2913	2917	2921	rVB2	501908	709797	8.85%	0.429%
85	20.286	2938	2941	2945	rBV	309278	367800	4.58%	0.222%
86	20.333	2946	2949	2953	rVB	319506	380128	4.74%	0.230%
87	20.739	3014	3018	3023	rBV2	644469	1035935	12.91%	0.626%
88	20.969	3051	3057	3066	rBV3	1848690	4601059	57.35%	2.780%
89	21.210	3095	3098	3103	rVB	493450	575548	7.17%	0.348%
90	21.292	3106	3112	3117	rVV3	3417304	8022102	100.00%	4.847%
91	21.339	3117	3120	3132	rVB	2550279	3987753	49.71%	2.410%
92	21.504	3143	3148	3153	rVB3	1362663	2114076	26.35%	1.277%
93	21.569	3154	3159	3164	rVV	2063971	2996322	37.35%	1.811%
94	21.622	3164	3168	3171	rVV	620928	938301	11.70%	0.567%
95	21.657	3171	3174	3183	rVB3	792019	1278950	15.94%	0.773%
96	22.045	3236	3240	3249	rBV3	499483	1065579	13.28%	0.644%
97	22.669	3342	3346	3353	rVB2	869535	1653730	20.61%	0.999%
98	23.310	3449	3455	3463	rBV	1865857	5279193	65.81%	3.190%
99	24.039	3573	3579	3585	rBV2	1641419	3455710	43.08%	2.088%
100	24.239	3607	3613	3640	rVB	1773363	5690216	70.93%	3.438%

Sum of corrected areas: 165490439

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
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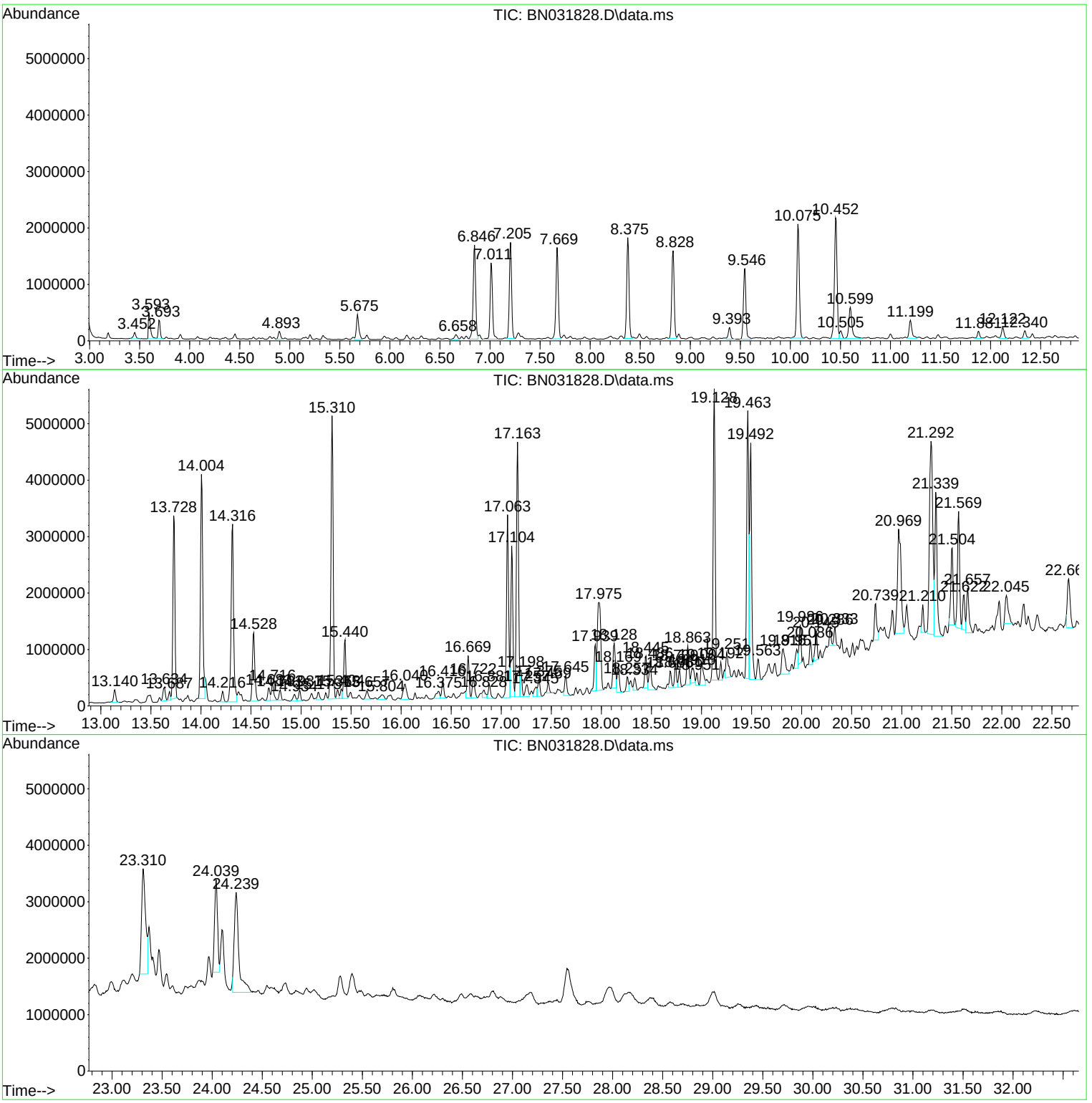
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

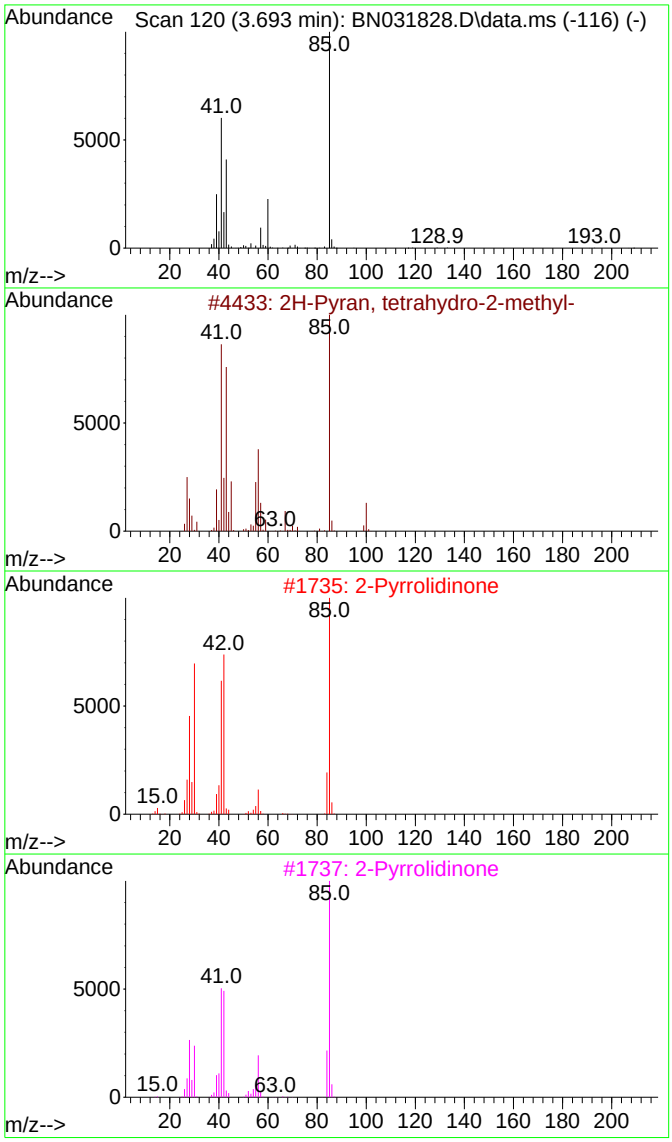
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

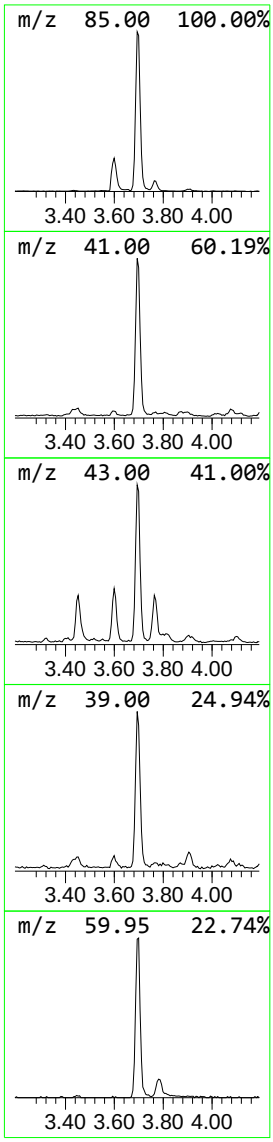
****

Peak Number 1 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.693	3.33 ng/ul	437123	1,4-Dichlorobenzene-d4	7.669

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
2	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
3	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
4	trans-4-Methoxy-3-buten-2-one	100	C5H8O2	051731-17-0	9
5	2(3H)-Furanone, 5-ethyldihydro-	114	C6H10O2	000695-06-7	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
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Sample : P2634-15
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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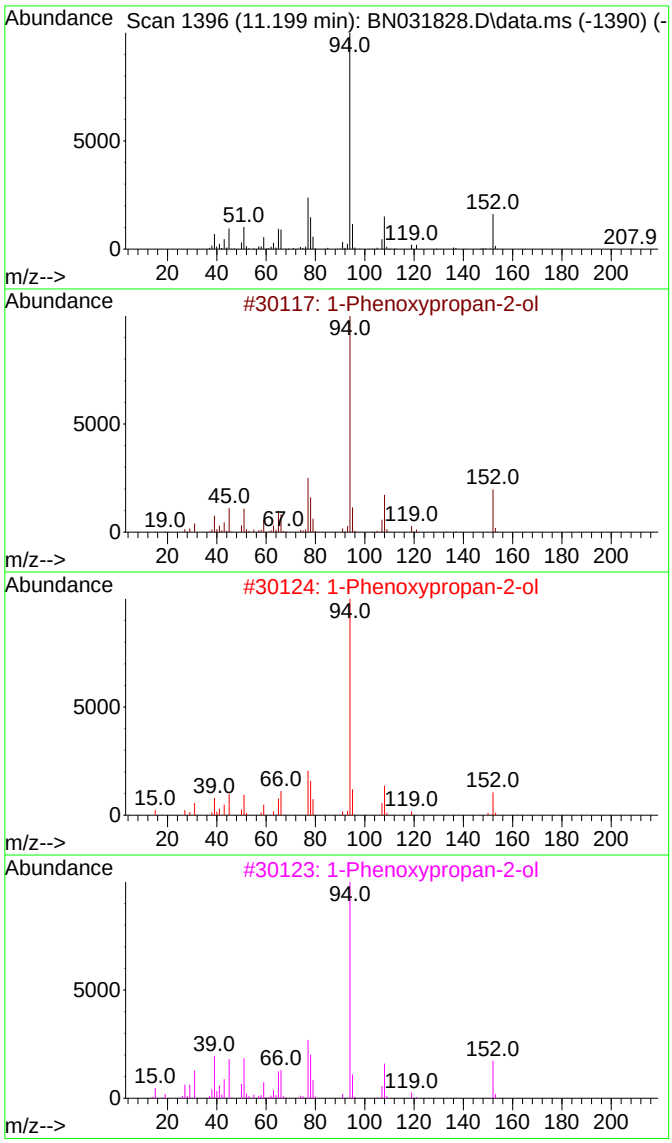
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

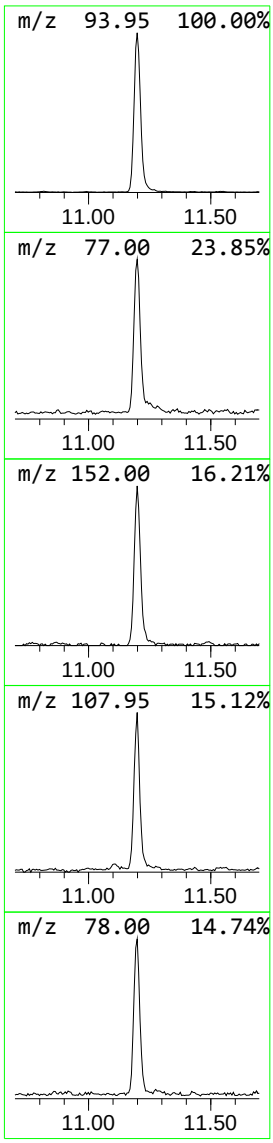
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.199	3.23 ng/ul	637721	Naphthalene-d8	10.458

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	95
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	90
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
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Sample : P2634-15
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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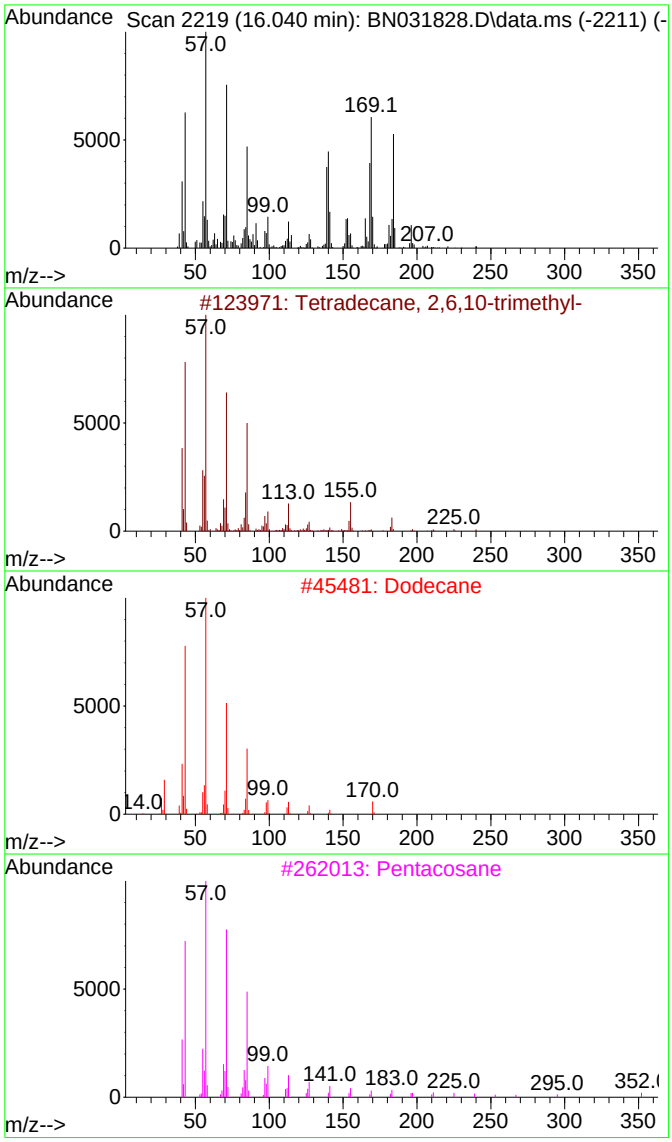
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

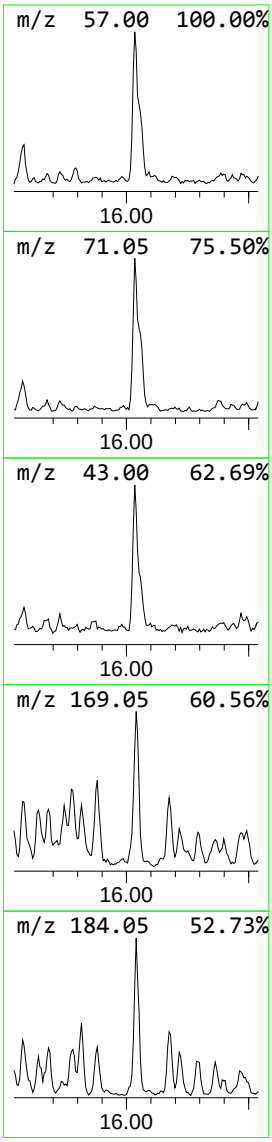
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.040	2.76 ng/ul	657487	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	52
2	Dodecane	170	C12H26	000112-40-3	42
3	Pentacosane	352	C25H52	000629-99-2	38
4	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	38
5	Dodecane	170	C12H26	000112-40-3	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
Acq On : 06 Jun 2024 21:12
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

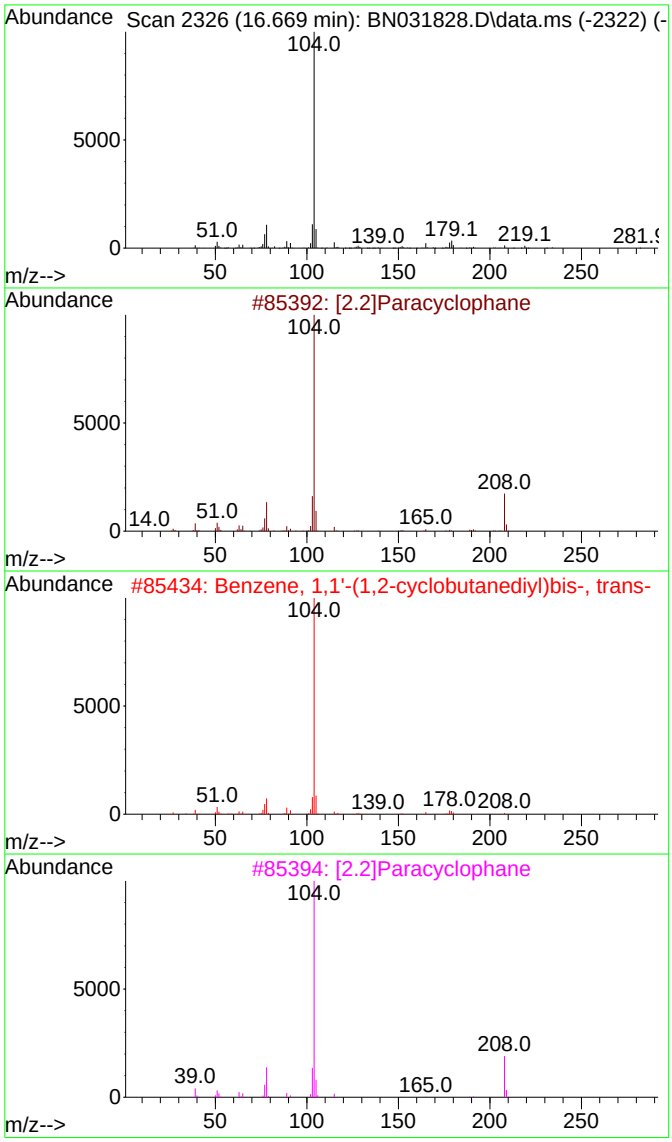
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

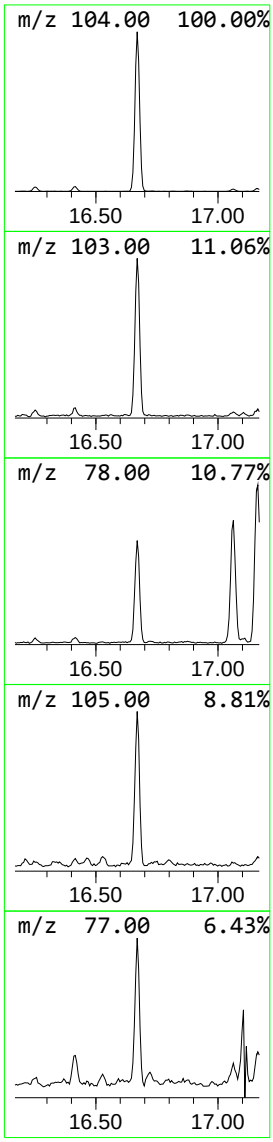
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 [2.2]Paracyclophane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	4.30 ng/ul	1026340	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[2.2]Paracyclophane	208	C16H16	001633-22-3	87
2	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	86
3	[2.2]Paracyclophane	208	C16H16	001633-22-3	83
4	Cyclobutane, 1,2-diphenyl-	208	C16H16	003018-21-1	72
5	[2.2]Paracyclophane	208	C16H16	001633-22-3	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

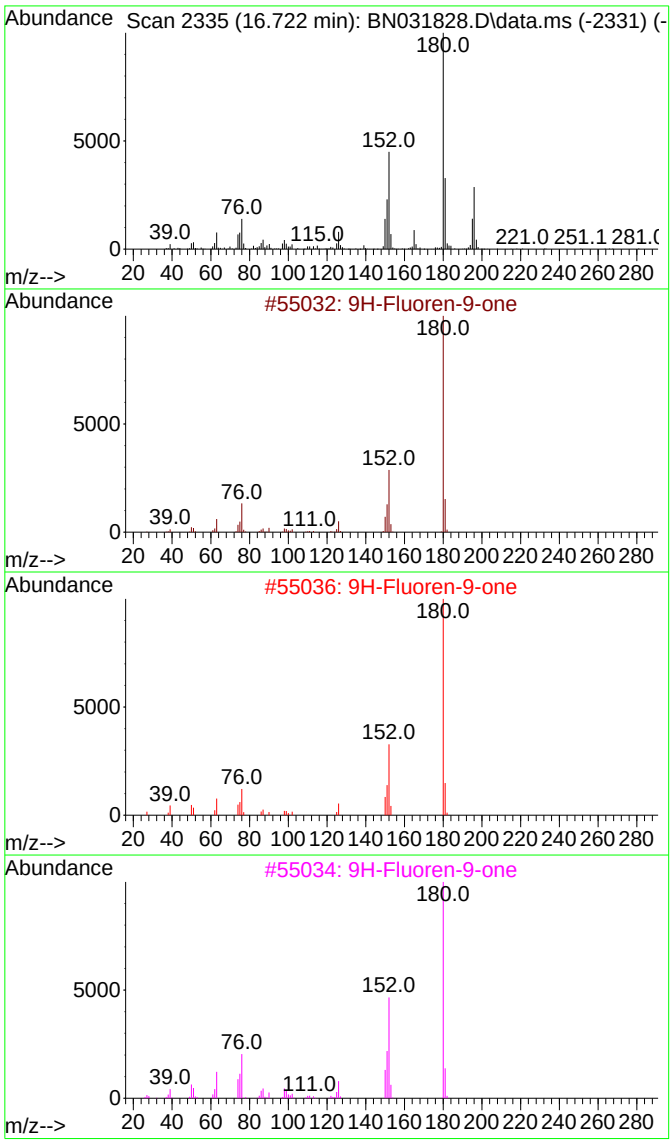
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

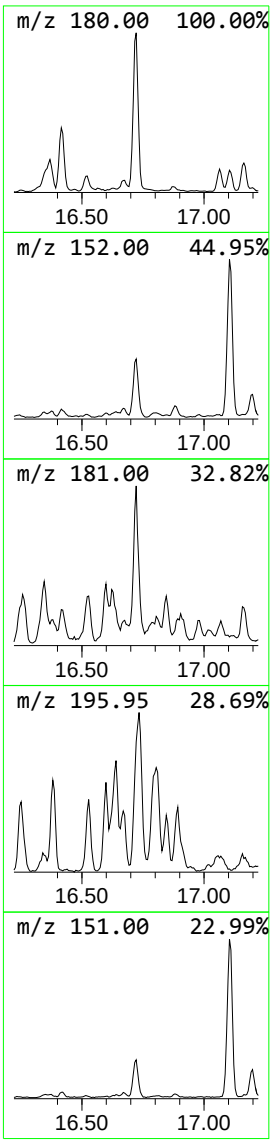
****

Peak Number 6 9H-Fluoren-9-one Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.722	2.49 ng/ul	593564	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	92
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	76
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	64
5	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	64





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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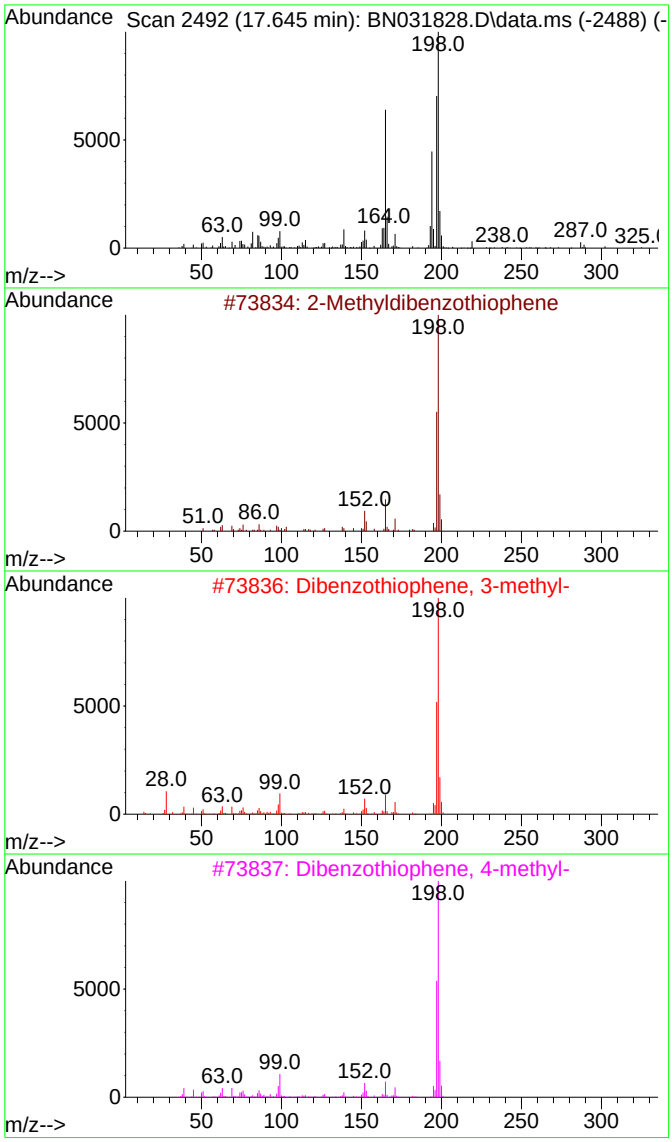
Quant Title : SVOA CALIBRATION

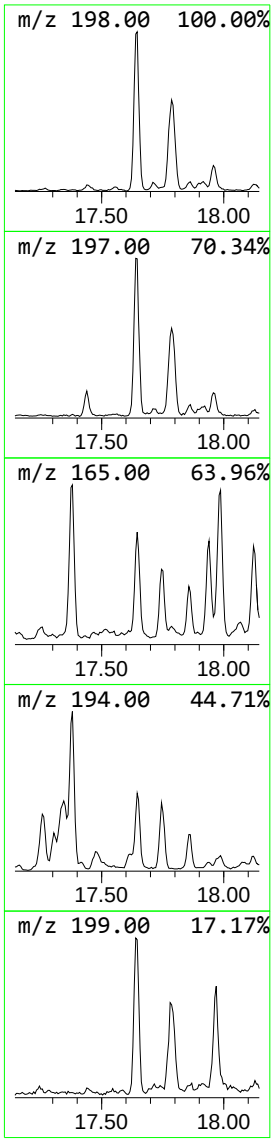
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Methyldibenzothiophene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.645	2.32 ng/ul	554563	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	62
2	Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	55
3	Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	55
4	Benzo[c]cinnoline, 4-methyl-	194	C13H10N2	019174-78-8	52
5	9H-Fluorene-9-thiol	198	C13H10S	019552-08-0	52





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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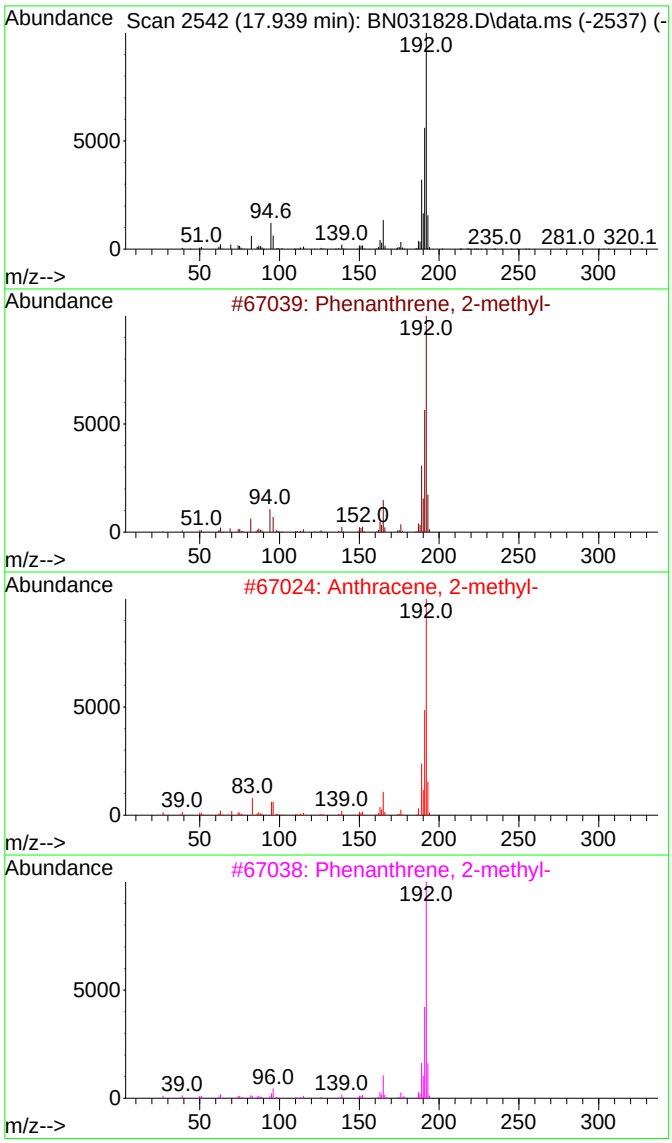
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

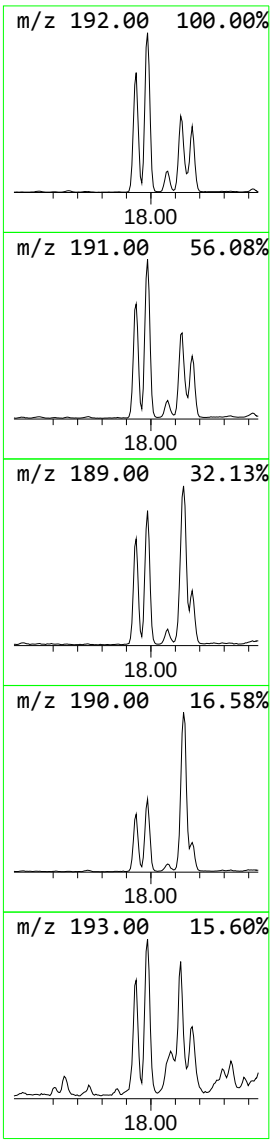
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.939	5.11 ng/ul	1218190	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93





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Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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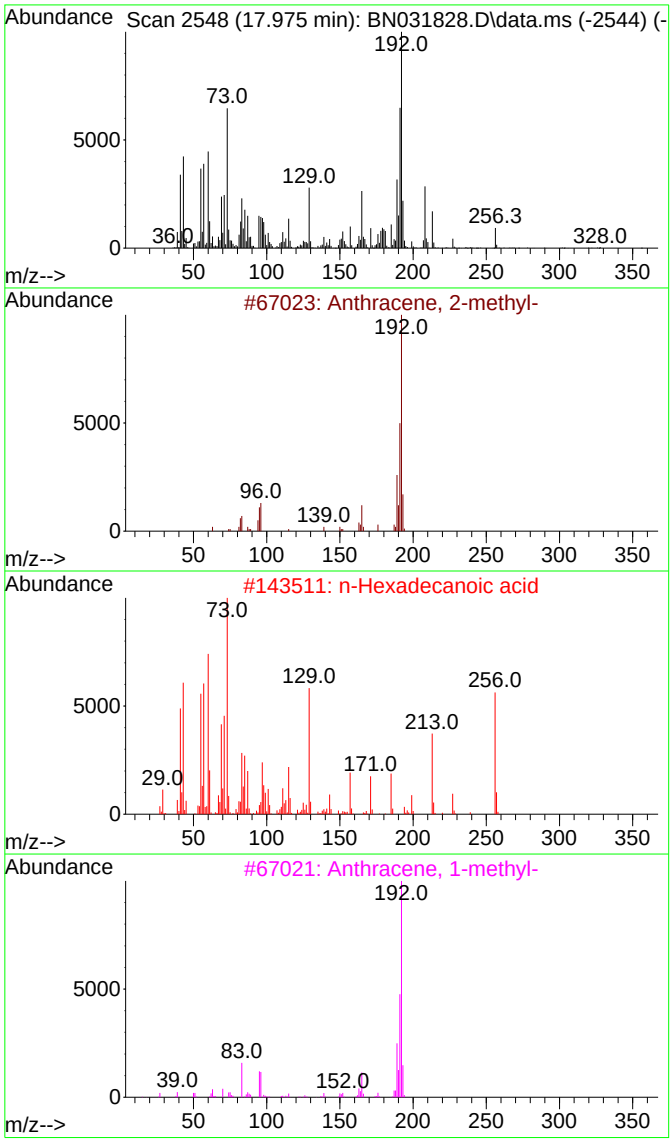
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

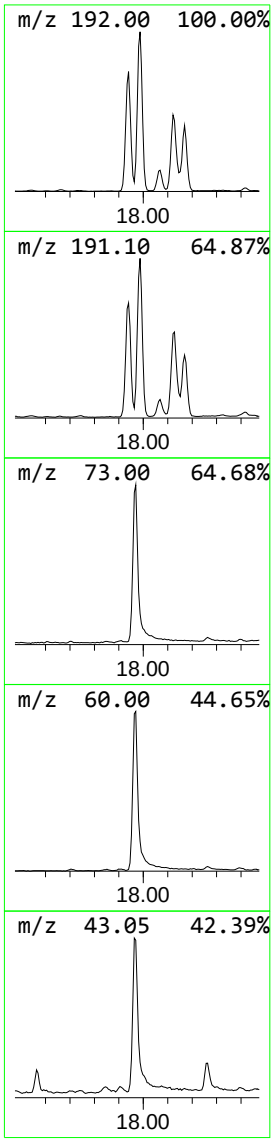
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Anthracene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.975	15.00 ng/ul	3579620	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	87
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	70





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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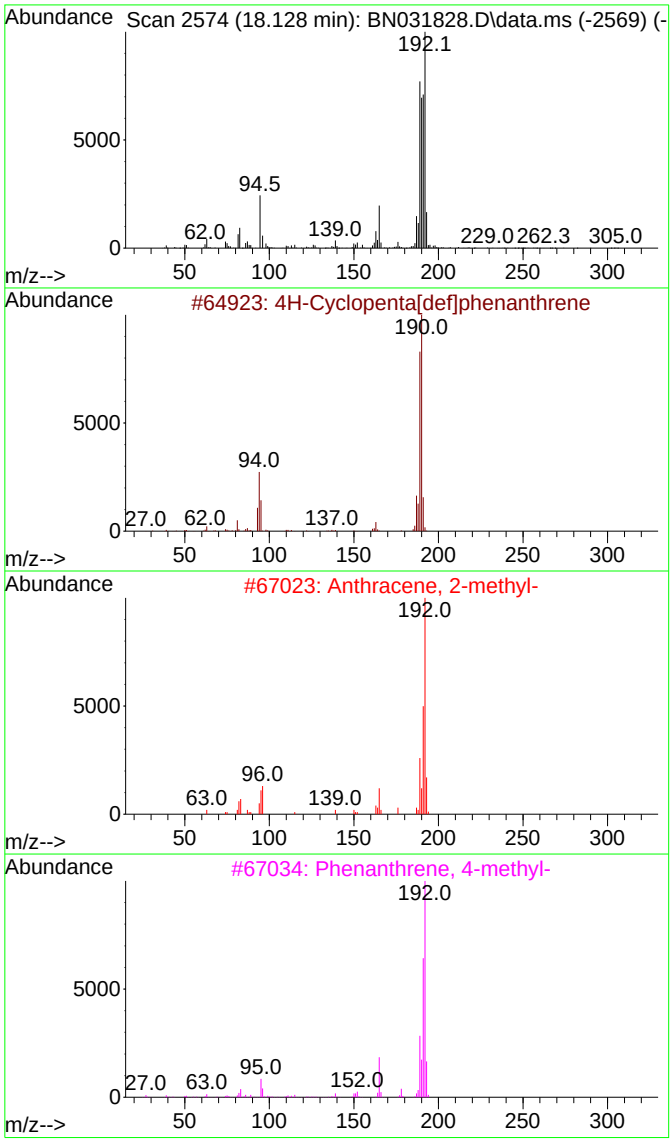
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

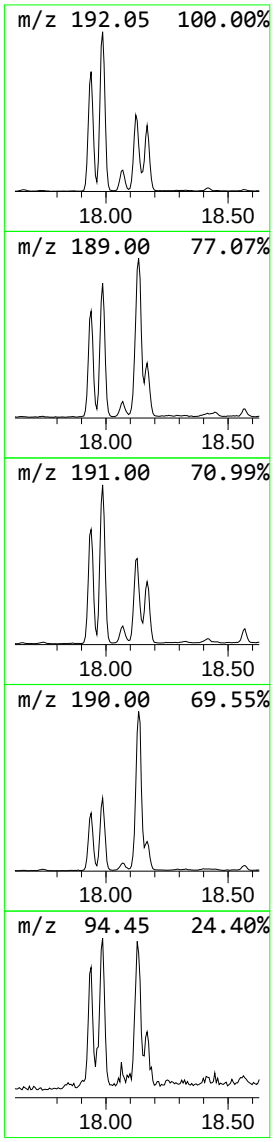
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	5.63 ng/ul	1343640	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	50
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	50
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	50
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	50





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Misc :
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ClientSampleId :
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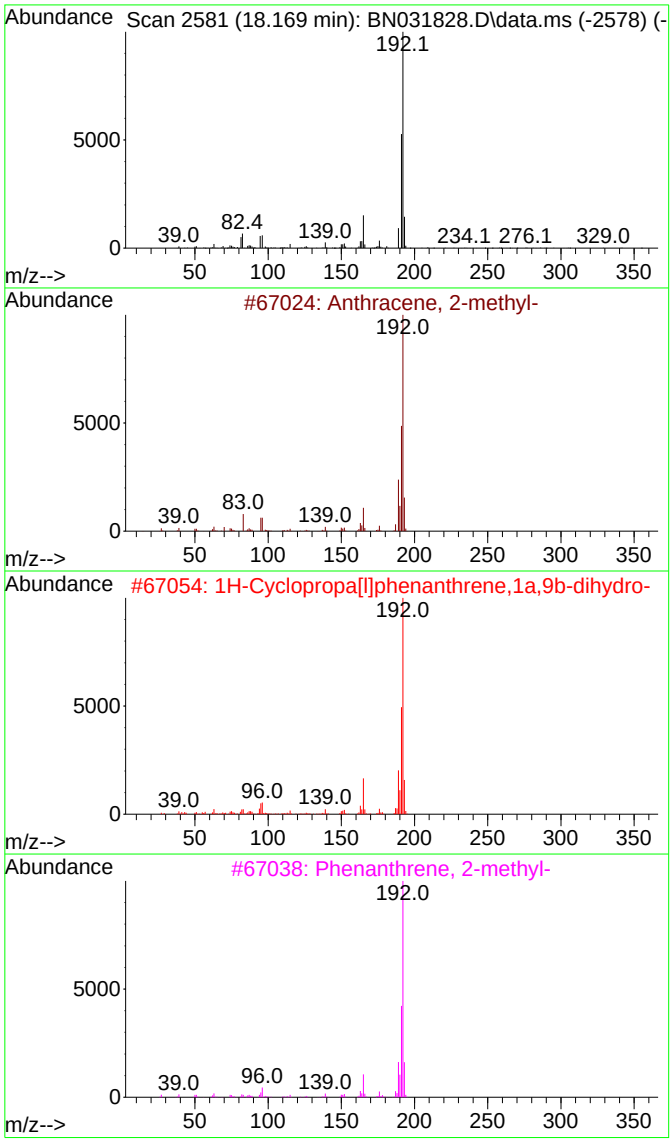
Quant Title : SVOA CALIBRATION

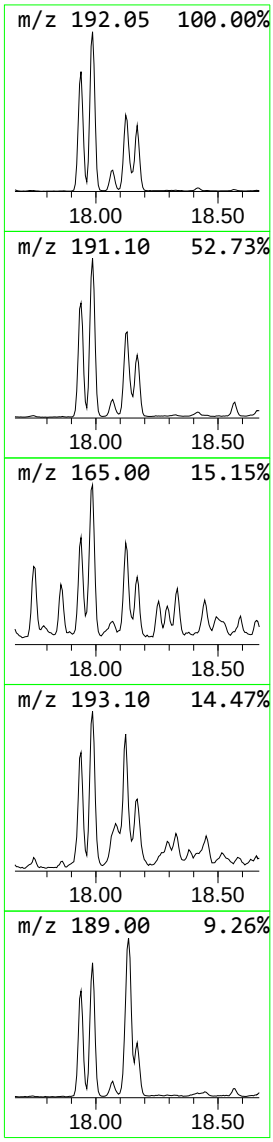
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1H-Cyclopropa[l]phenanthren... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	2.80 ng/ul	667571	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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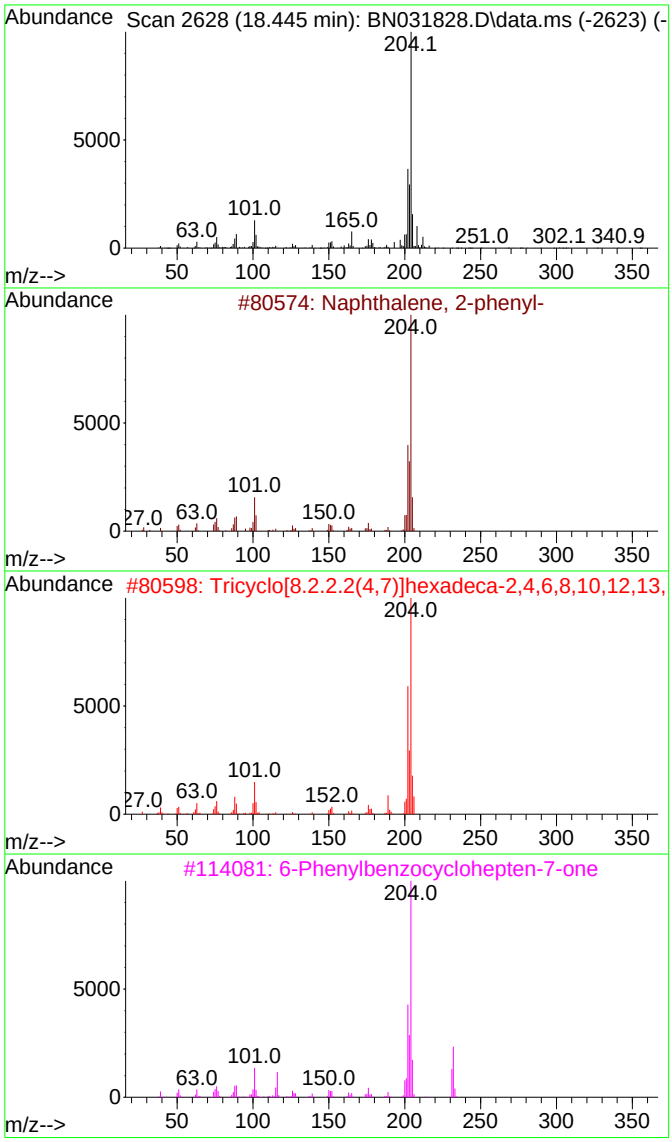
Quant Title : SVOA CALIBRATION

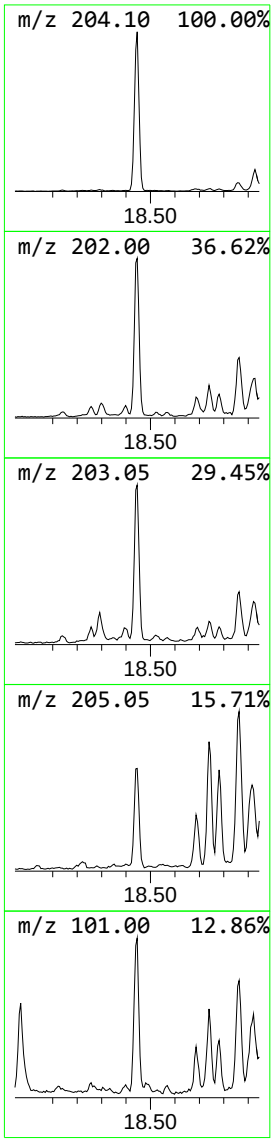
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.445	3.39 ng/ul	809068	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	89
3	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86
4	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

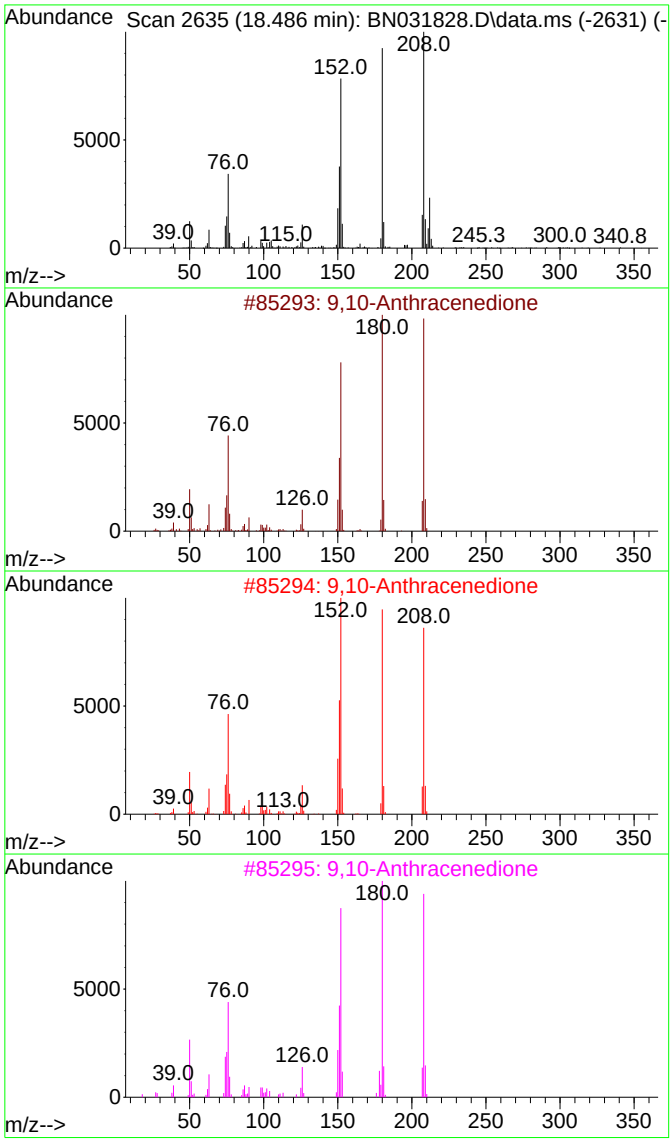
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

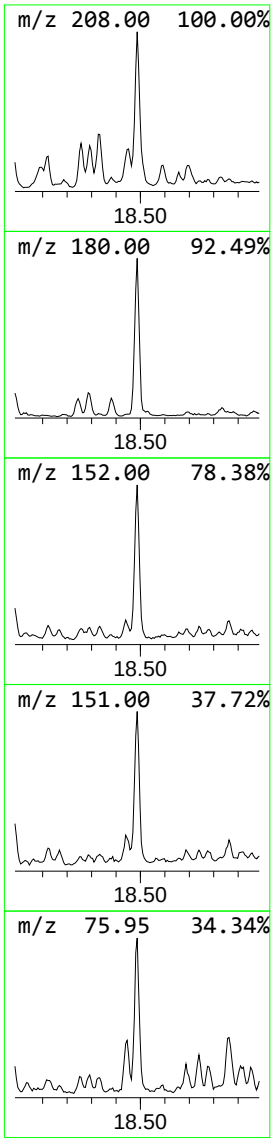
****

Peak Number 13 9,10-Anthracenedione Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	3.51 ng/ul	838037	Phenanthrene-d10	17.063

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	98
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	55





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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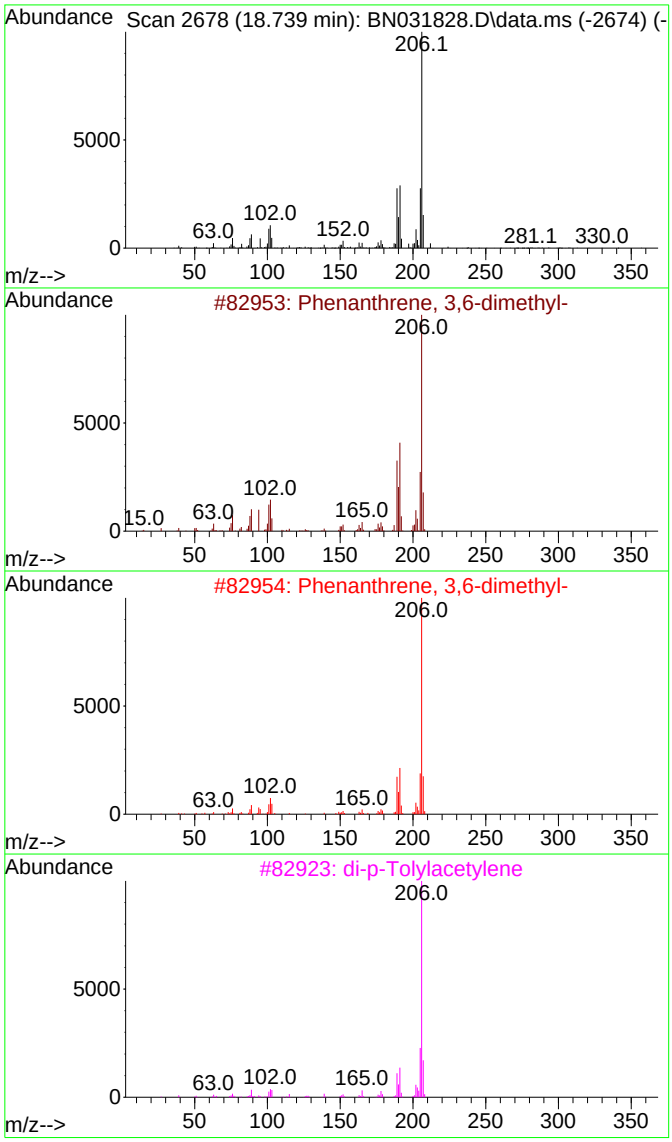
Quant Title : SVOA CALIBRATION

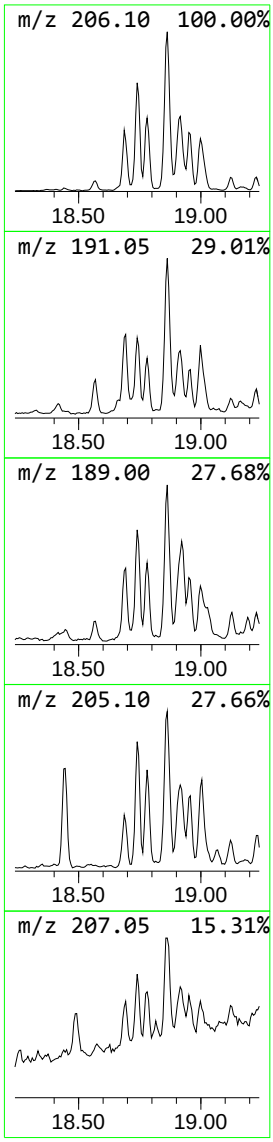
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Phenanthrene, 3,6-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.739	2.38 ng/ul	568619	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
4	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	87
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	83





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ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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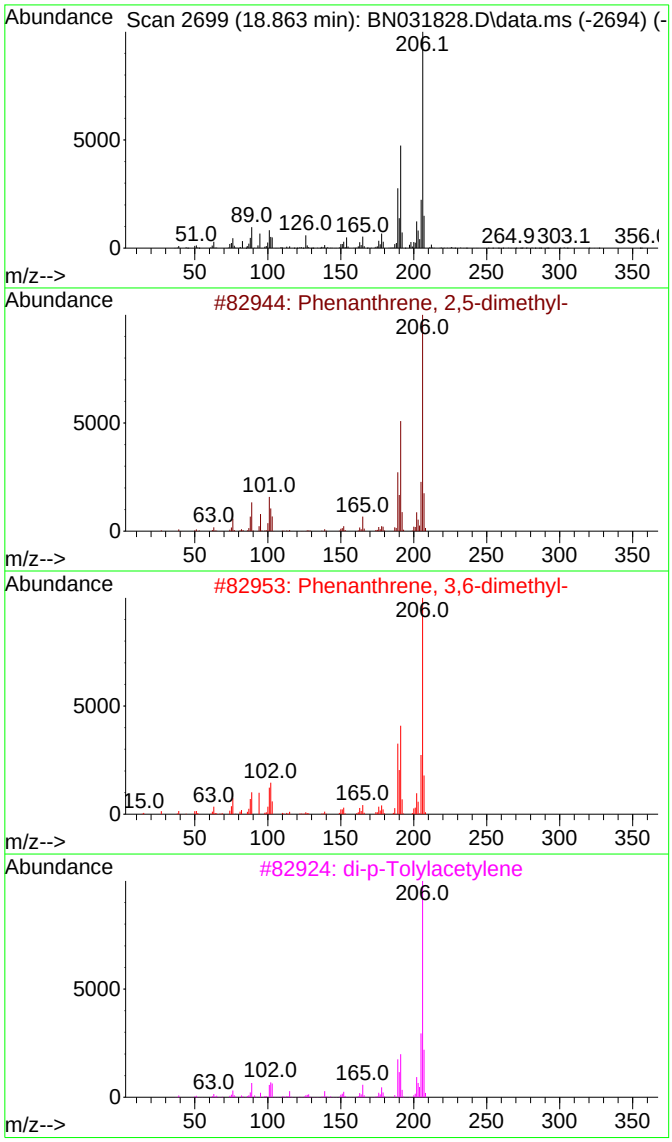
Quant Title : SVOA CALIBRATION

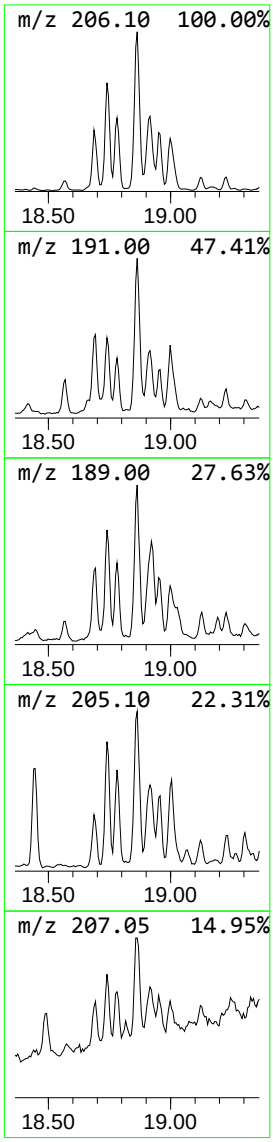
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.863	4.44 ng/ul	1058330	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
5	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90





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Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

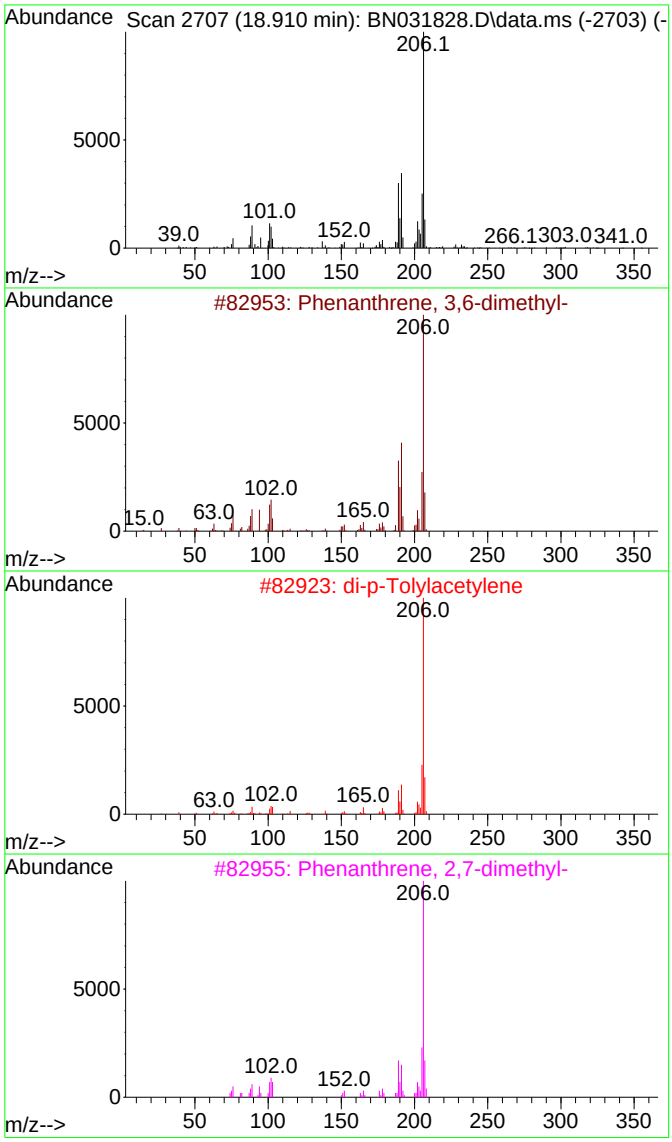
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

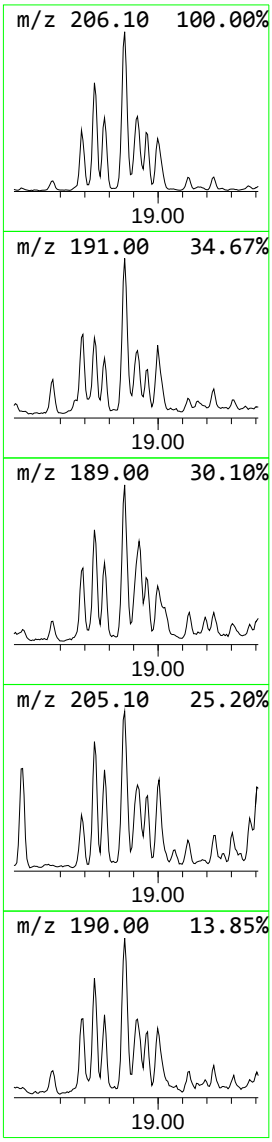
****

Peak Number 16 di-p-Tolylacetylene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.910	2.32 ng/ul	554161	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	96
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
3	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	90
4	di-p-Tolylacetylene	206	C16H14	002789-88-0	90
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	83





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ALS Vial : 16 Sample Multiplier: 1

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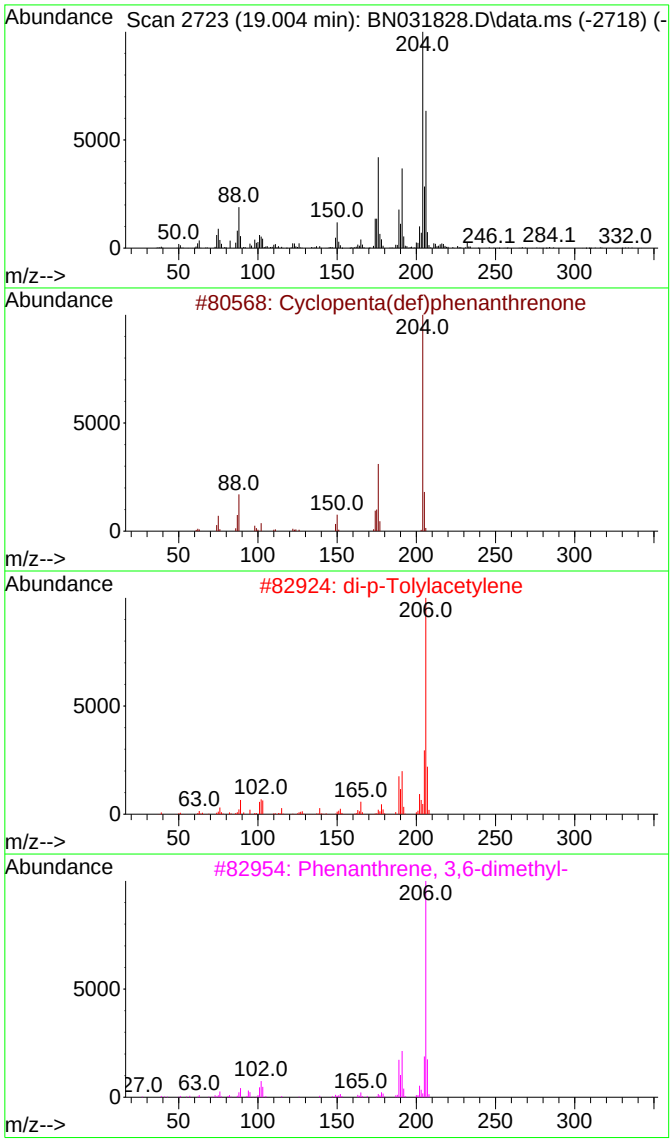
Quant Title : SVOA CALIBRATION

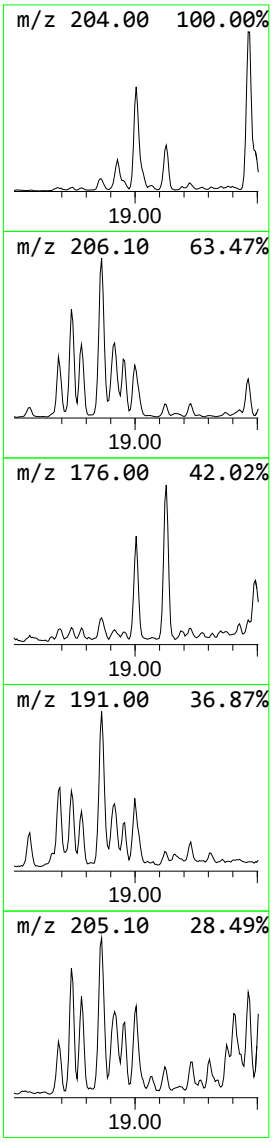
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Cyclopenta(def)phenanthrenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	3.28 ng/ul	782917	Phenanthrene-d10	17.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	95
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	38
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	38
4	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	30
5	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	30





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Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
BHBX9

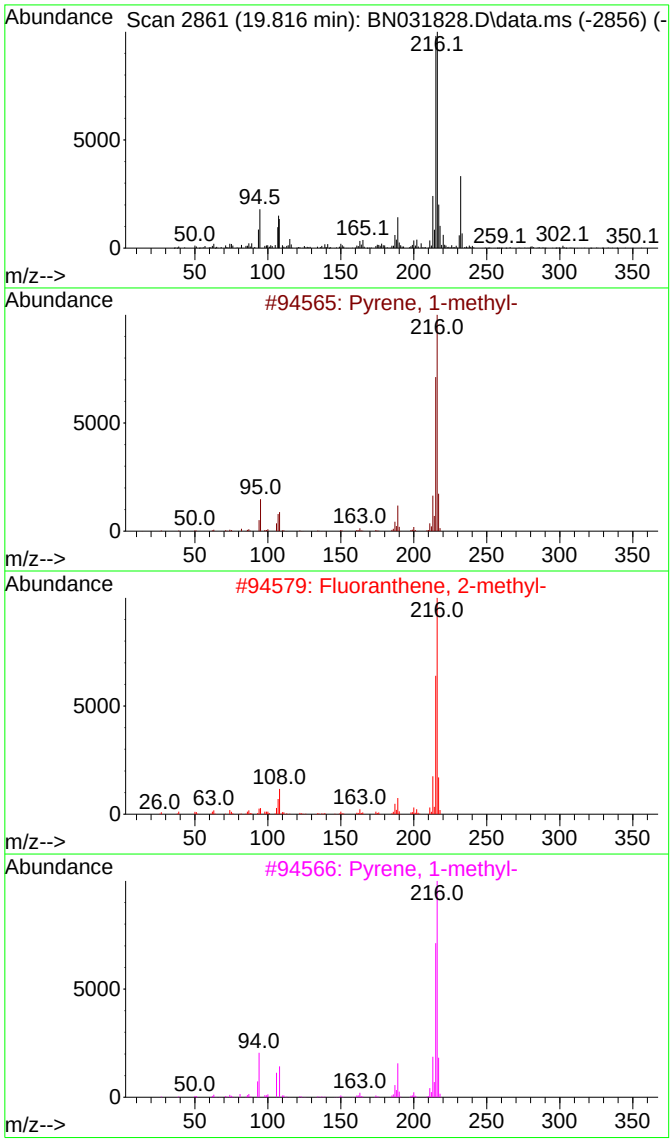
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

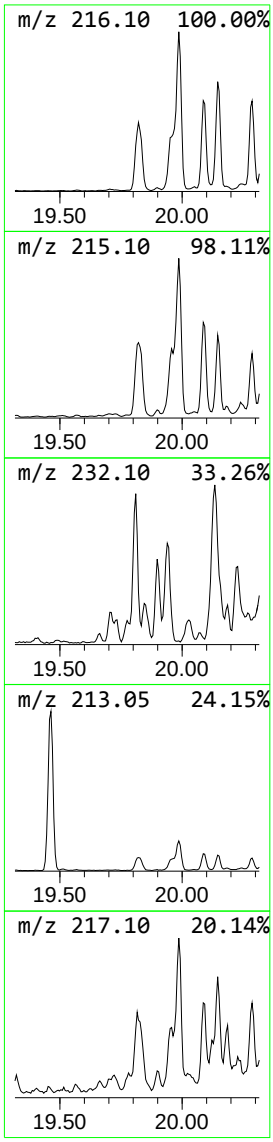
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Pyrene, 1-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.816	2.76 ng/ul	1108260	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	89
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

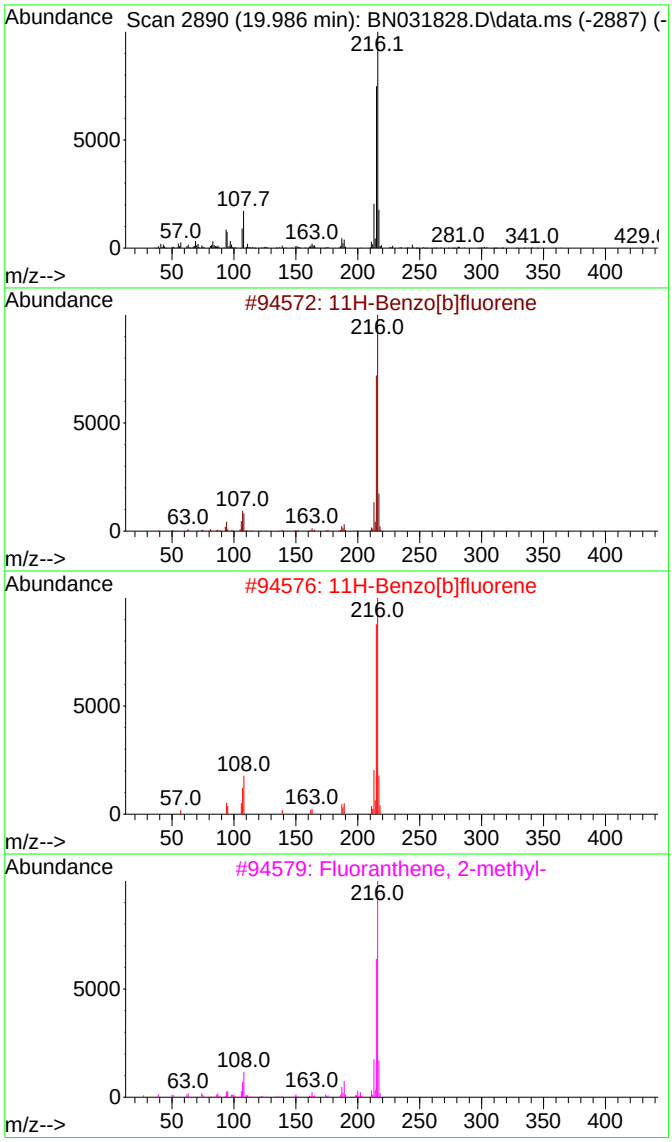
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

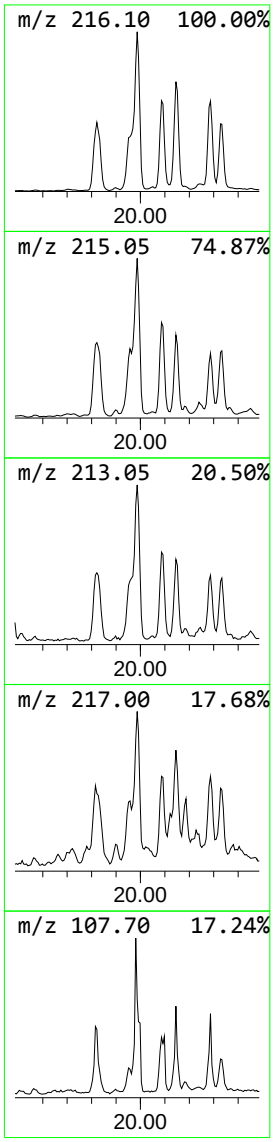
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 11H-Benzo[b]fluorene Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.986	2.06 ng/ul	827588	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

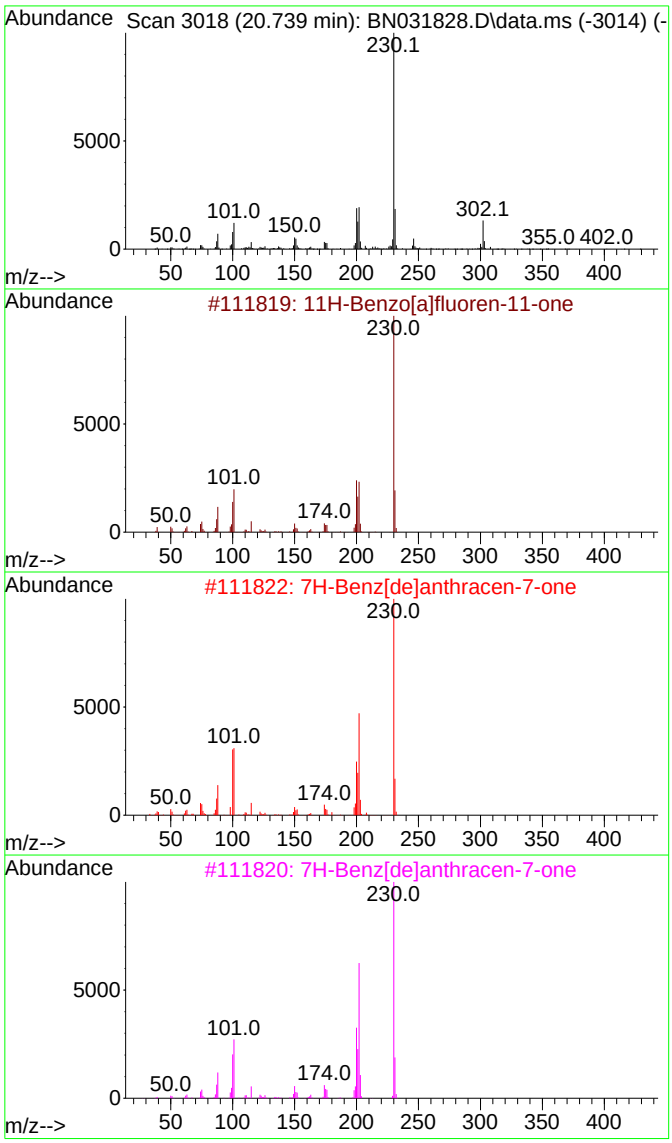
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

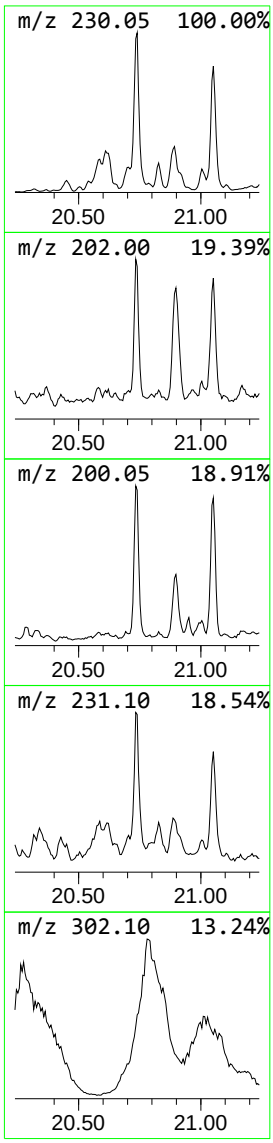
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 11H-Benzo[a]fluoren-11-one Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	2.58 ng/ul	1035940	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	98
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	86
5	4-Pyrenecarbaldehyde	230	C17H10O	022245-51-8	72





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Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

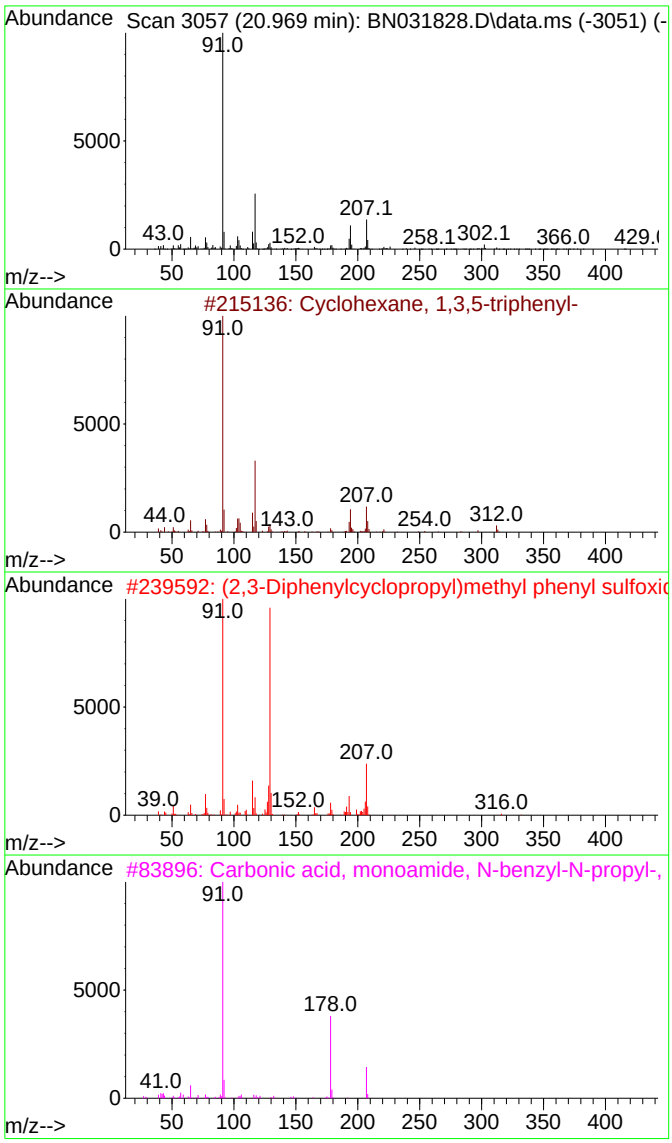
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

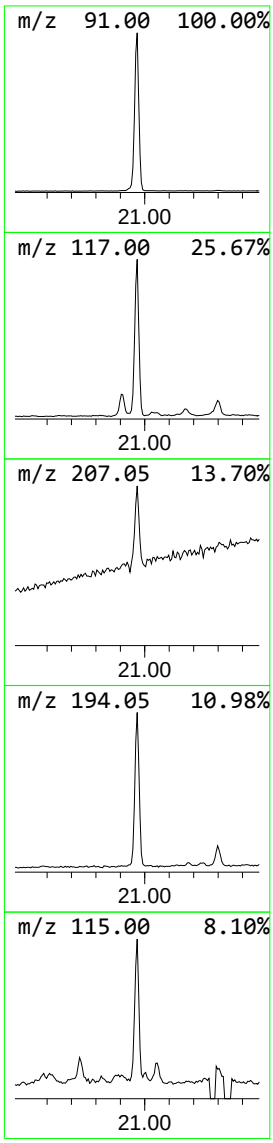
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Cyclohexane, 1,3,5-triphenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.969	11.47 ng/ul	4601060	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3,5-triphenyl-	312	C24H24	028336-57-4	83
2	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	38
3	Carbonic acid, monoamide, N-benz...	207	C12H17NO2	1000451-48-5	27
4	(R)-(1-Cyclopentylpropane-1,3-di...	264	C20H24	1402851-74-4	16
5	1-benzylindole	207	C15H13N	1000334-31-3	16





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

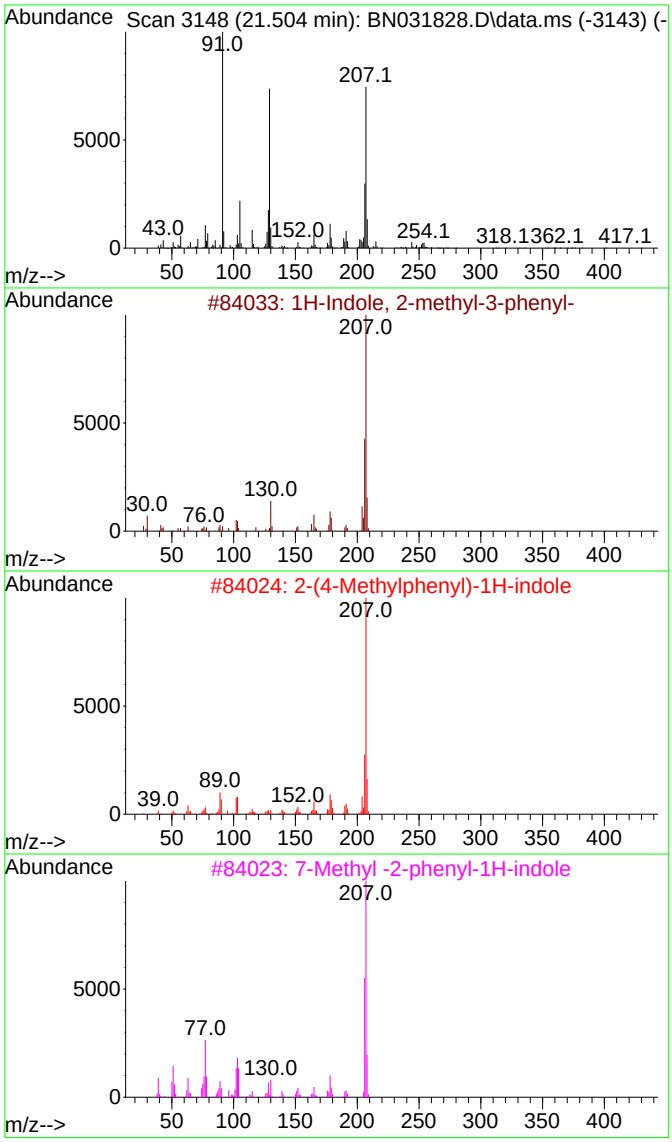
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

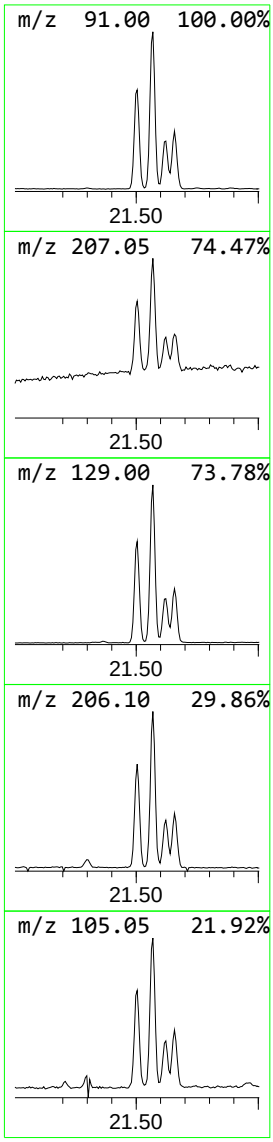
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 1H-Indole, 2-methyl-3-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.504	5.27 ng/ul	2114080	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	50
2	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	46
3	7-Methyl -2-phenyl-1H-indole	207	C15H13N	059541-82-1	43
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	43
5	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	42





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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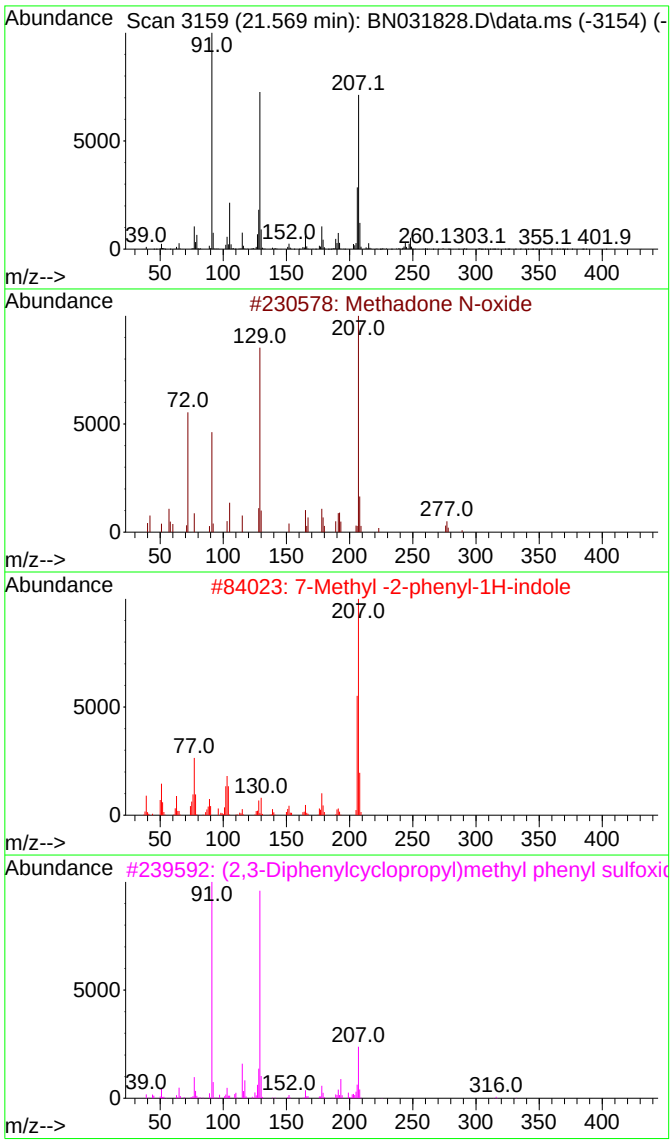
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

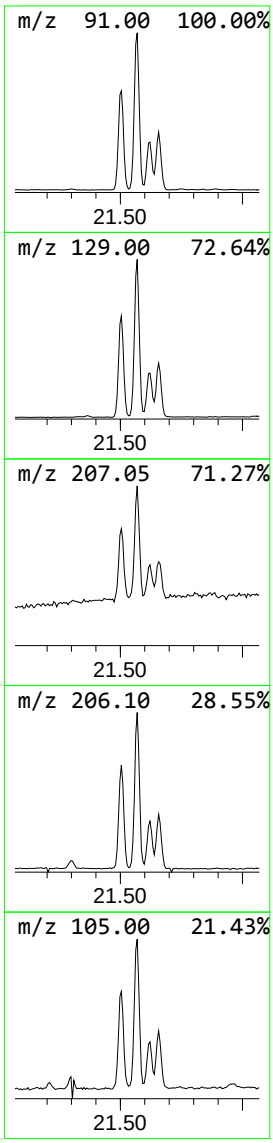
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Methadone N-oxide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.569	7.47 ng/ul	2996320	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	52
2	7-Methyl -2-phenyl-1H-indole	207	C15H13N	059541-82-1	43
3	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	42
4	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	41
5	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	41





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

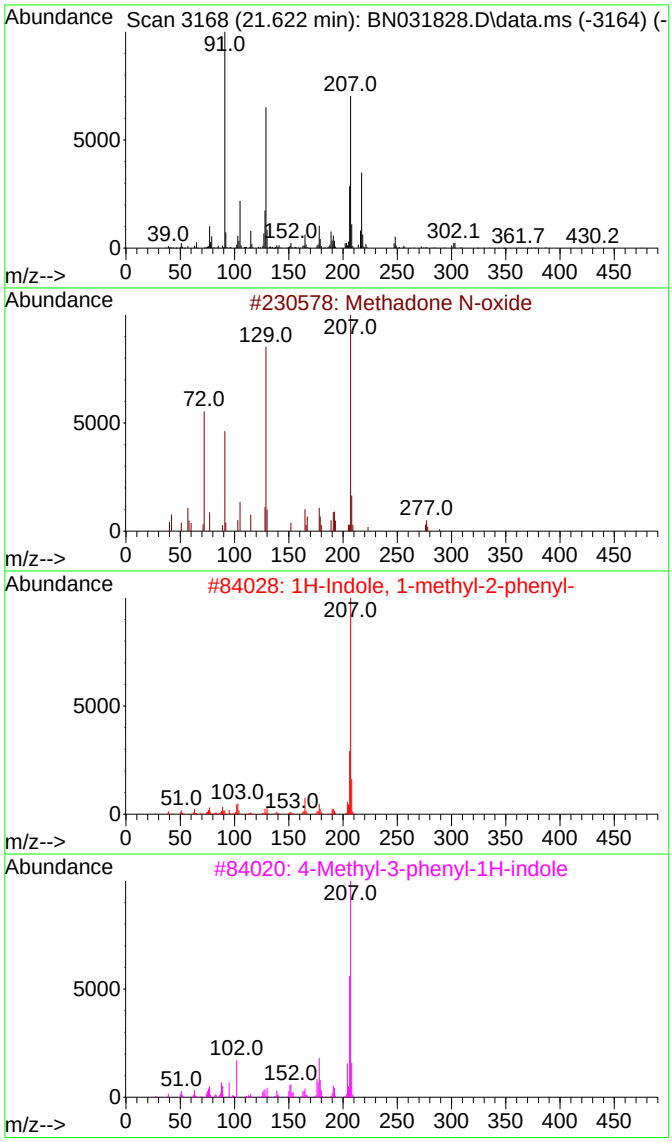
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

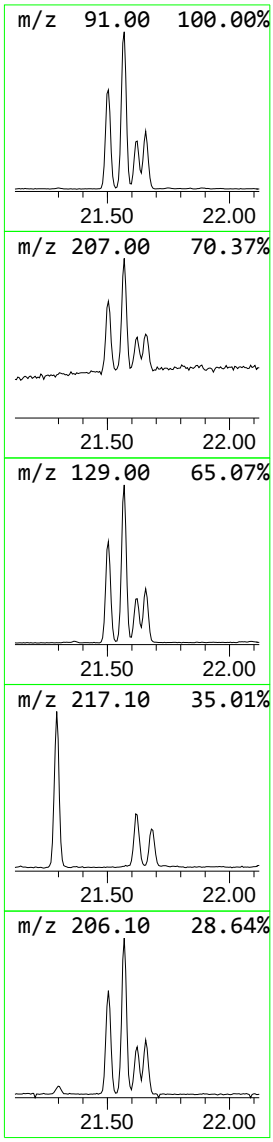
****

Peak Number 24 unknown-02 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.622	2.34 ng/ul	938301	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	47
2	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	42
3	4-Methyl-3-phenyl-1H-indole	207	C15H13N	180271-54-9	38
4	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	38
5	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	35





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

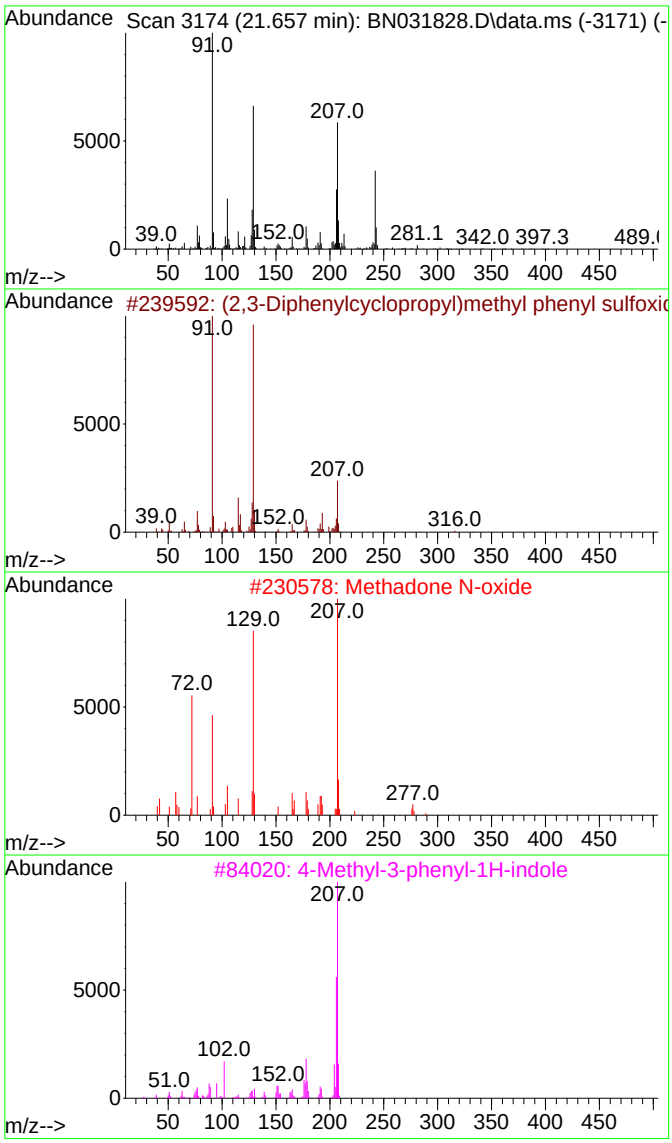
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

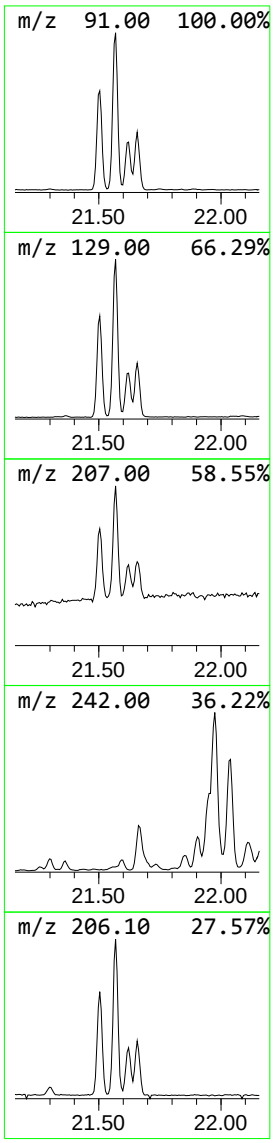
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (2,3-Diphenylcyclopropyl)me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.657	3.19 ng/ul	1278950	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	50
2	Methadone N-oxide	325	C21H27NO2	1000120-80-7	46
3	4-Methyl-3-phenyl-1H-indole	207	C15H13N	180271-54-9	27
4	3-Hydroxytriazolo[4,3-a]pyridine...	207	C9H13N3OSi	1000476-39-3	25
5	1H-Indole, 5-methyl-2-phenyl-	207	C15H13N	013228-36-9	25





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

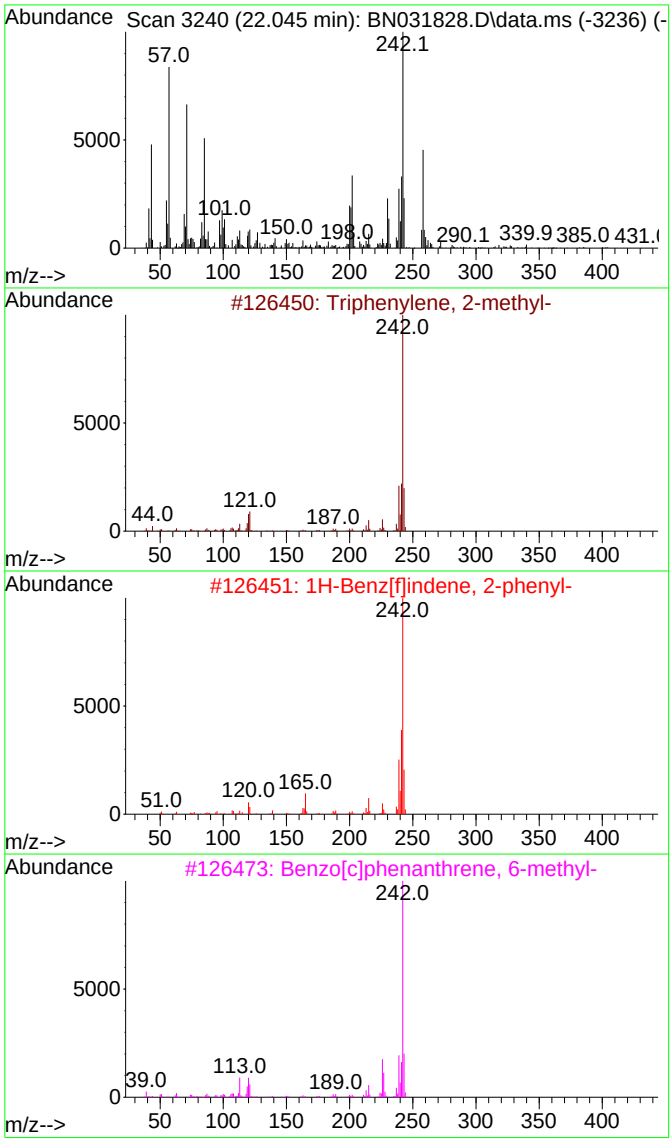
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

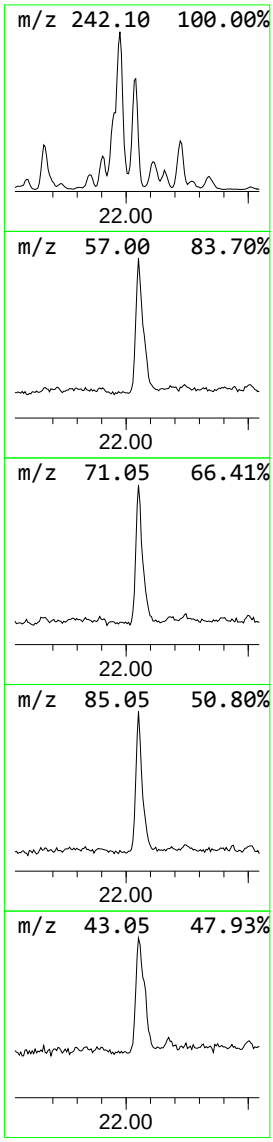
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Triphenylene, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	2.66 ng/ul	1065580	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93
2	1H-Benz[f]indene, 2-phenyl-	242	C19H14	1000305-22-4	86
3	Benzo[c]phenanthrene, 6-methyl-	242	C19H14	002381-34-2	80
4	Benz[a]anthracene, 12-methyl-	242	C19H14	002422-79-9	70
5	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	66





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Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

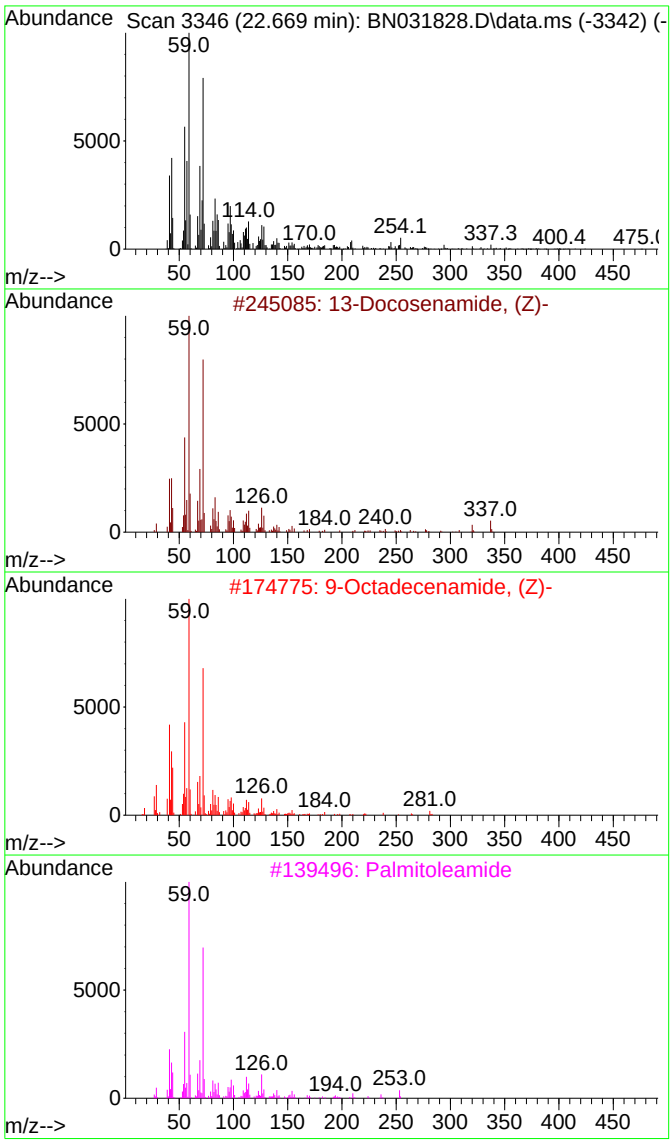
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

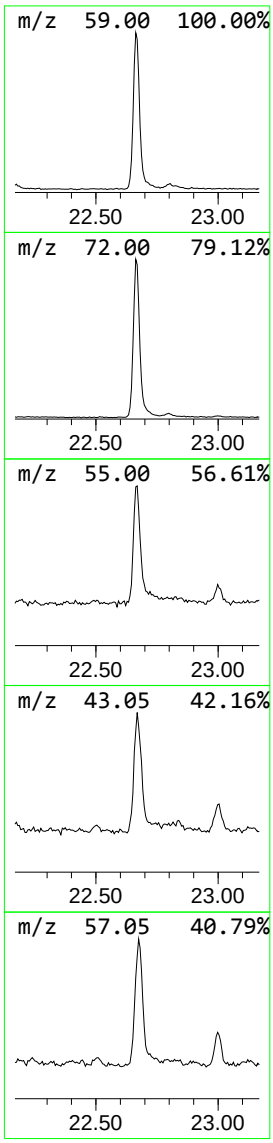
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 13-Docosenamide, (Z)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.669	4.12 ng/ul	1653730	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	97
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	83
3	Palmitoleamide	253	C16H31NO	106010-22-4	80
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72
5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031828.D
Acq On : 06 Jun 2024 21:12
Operator : MA/JU
Sample : P2634-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	3.693	3.3	ng/ul	437123	1	7.669	2622650	20.0
1-Phenoxypropan...	11.199	3.2	ng/ul	637721	2	10.458	3944440	20.0
(DEL) Alkane: S...	16.040	2.8	ng/ul	657487	4	17.063	4772130	20.0
[2.2]Paracyclop...	16.669	4.3	ng/ul	1026340	4	17.063	4772130	20.0
9H-Fluoren-9-one	16.722	2.5	ng/ul	593564	4	17.063	4772130	20.0
2-Methyldibenzo...	17.645	2.3	ng/ul	554563	4	17.063	4772130	20.0
Phenanthrene, 2...	17.939	5.1	ng/ul	1218190	4	17.063	4772130	20.0
Anthracene, 2-m...	17.975	15.0	ng/ul	3579620	4	17.063	4772130	20.0
4H-Cyclopenta[d...	18.128	5.6	ng/ul	1343640	4	17.063	4772130	20.0
1H-Cyclopropa[l...	18.169	2.8	ng/ul	667571	4	17.063	4772130	20.0
Naphthalene, 2-...	18.445	3.4	ng/ul	809068	4	17.063	4772130	20.0
9,10-Anthracene...	18.486	3.5	ng/ul	838037	4	17.063	4772130	20.0
Phenanthrene, 3...	18.739	2.4	ng/ul	568619	4	17.063	4772130	20.0
Phenanthrene, 2...	18.863	4.4	ng/ul	1058330	4	17.063	4772130	20.0
di-p-Tolylacety...	18.910	2.3	ng/ul	554161	4	17.063	4772130	20.0
Cyclopenta(def)...	19.004	3.3	ng/ul	782917	4	17.063	4772130	20.0
Pyrene, 1-methyl-	19.816	2.8	ng/ul	1108260	5	21.298	8022100	20.0
11H-Benzo[b]flu...	19.986	2.1	ng/ul	827588	5	21.298	8022100	20.0
11H-Benzo[a]flu...	20.739	2.6	ng/ul	1035940	5	21.298	8022100	20.0
Cyclohexane, 1,...	20.969	11.5	ng/ul	4601060	5	21.298	8022100	20.0
1H-Indole, 2-me...	21.504	5.3	ng/ul	2114080	5	21.298	8022100	20.0
Methadone N-oxide	21.569	7.5	ng/ul	2996320	5	21.298	8022100	20.0
unknown-02	21.622	2.3	ng/ul	938301	5	21.298	8022100	20.0
(2,3-Diphenylcy...	21.657	3.2	ng/ul	1278950	5	21.298	8022100	20.0
Triphenylene, 2...	22.045	2.7	ng/ul	1065580	5	21.298	8022100	20.0
13-Docosenamide...	22.669	4.1	ng/ul	1653730	5	21.298	8022100	20.0