LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM047059.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	19	22	27	rBV	55436	69042	0.86%	0.068%
2	3.416	86	90	104	rBV	247049	406701	5.07%	0.400%
3	6.598	620	631	641	rBV	527679	813241	10.14%	0.799%
4	6.751	648	657	668	rBV	993739	1488437	18.56%	1.463%
5	6.934	681	688	702	rBV	1224370	1900718	23.70%	1.868%
6	7.393	759	766	774	rBV	1107470	1672790	20.86%	1.644%
7	7.757	822	828	835	rVB	154863	244515	3.05%	0.240%
8	8.116	881	889	898	rBV	1201156	1886522	23.52%	1.854%
9	8.539	954	961	970	rVB	1150256	1856242	23.14%	1.824%
10	8.728	988	993	1000	rVB2	43731	68904	0.86%	0.068%
11	8.857	1004	1015	1020	rVB3	15298	39832	0.50%	0.039%
12	9.122	1055	1060	1065	rBV2	41876	64434	0.80%	0.063%
13	9.251	1075	1082	1089	rBV	987087	1599714	19.95%	1.572%
14	9.569	1130	1136	1145	rVB5	34345	75326	0.94%	0.074%
15	9.786	1166	1173	1184	rVV	1565449	2534131	31.60%	2.490%
16	10.151	1226	1235	1240	rBV	1407366	2365976	29.50%	2.325%
17	10.198	1240	1243	1251	rVB	116588	191251	2.38%	0.188%
18	10.298	1251	1260	1275	rBV	865061	1650930	20.59%	1.622%
19	10.404	1275	1278	1282	rVV2	37385	59270	0.74%	0.058%
20	10.451	1282	1286	1293	rVV3	20997	40958	0.51%	0.040%
21	10.569	1300	1306	1312	rVV3	31623	68644	0.86%	0.067%
22	10.728	1327	1333	1339	rVB5	25418	47648	0.59%	0.047%
23	10.798	1339	1345	1353	rVB4	15381	38241	0.48%	0.038%
24	10.910	1357	1364	1372	rBV	74312	138473	1.73%	0.136%
25	11.251	1416	1422	1428	rBV	97987	163264	2.04%	0.160%
26	11.651	1482	1490	1495	rBV9	16146	40140	0.50%	0.039%
27	11.716	1495	1501	1505	rBV	63816	120074	1.50%	0.118%
28	12.051	1547	1558	1568	rVB	421340	834637	10.41%	0.820%
29	12.280	1594	1597	1605	rVB4	23296	53227	0.66%	0.052%
30	12.380	1608	1614	1619	rBV9	18009	39418	0.49%	0.039%
31	12.874	1693	1698	1704	rVV9	23157	39479	0.49%	0.039%
32	13.098	1731	1736	1745	rVB3	117778	220040	2.74%	0.216%
33	13.210	1745	1755	1758	rBV6	49053	116610	1.45%	0.115%
34	13.363	1775	1781	1787	rBV4	37863	87000	1.08%	0.085%
35	13.410	1787	1789	1792	rVV3	24840	37379	0.47%	0.037%
36	13.474	1792	1800	1807	rVB	2530760	3633088	45.30%	3.570%

LSC Area Percent Report

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Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Title : SVOA CALIBRATION

37	13.598	1817	1821	1824	rBV	82535	119745	1.49%	0.118%
38	13.633	1824	1827	1833	rVB	80277	116117	1.45%	0.114%
39	13.733	1837	1844	1859	rBV	2829277	4444847	55.42%	4.368%
40	13.998	1882	1889	1890	rBV6	21235	37445	0.47%	0.037%
41	14.051	1890	1898	1903	rVV	2163758	3346016	41.72%	3.288%
42	14.116	1903	1909	1915	rVV	5548058	8020057	100.00%	7.880%
43	14.186	1918	1921	1926	rVB2	54115	83834	1.05%	0.082%
44	14.292	1934	1939	1946	rBV2	431081	698765	8.71%	0.687%
45	14.369	1947	1952	1957	rVB	188047	280430	3.50%	0.276%
46	14.451	1959	1966	1974	rBV	946980	1410911	17.59%	1.386%
47	14.586	1980	1989	1997	rVB2	137004	329813	4.11%	0.324%
48	14.739	2011	2015	2017	rBV2	39680	59140	0.74%	0.058%
49	14.774	2017	2021	2026	rVB	133735	191784	2.39%	0.188%
50	15.016	2057	2062	2063	rBV	37874	59190	0.74%	0.058%
51	15.057	2063	2069	2073	rVV	3643832	5267926	65.68%	5.176%
52	15.110	2073	2078	2084	rVB	2640050	3626189	45.21%	3.563%
53	15.192	2086	2092	2097	rBV	1264175	1768369	22.05%	1.738%
54	15.233	2097	2099	2102	rVV2	155111	172675	2.15%	0.170%
55	15.268	2102	2105	2111	rVB	187599	238608	2.98%	0.234%
56	15.321	2111	2114	2118	rBV5	43347	76836	0.96%	0.075%
57	15.410	2124	2129	2134	rBV	439913	592288	7.39%	0.582%
58	15.516	2142	2147	2149	rBV3	39916	65126	0.81%	0.064%
59	15.563	2149	2155	2161	rVB3	302554	633409	7.90%	0.622%
60	15.621	2161	2165	2168	rBV	206321	315941	3.94%	0.310%
61	15.704	2175	2179	2183	rBV3	48098	65341	0.81%	0.064%
62	15.786	2188	2193	2203	rVB2	467578	772238	9.63%	0.759%
63	15.910	2208	2214	2222	rVB	470815	688376	8.58%	0.676%
64	16.074	2238	2242	2245	rBV3	66758	106459	1.33%	0.105%
65	16.121	2247	2250	2255	rVV2	81268	121613	1.52%	0.119%
66	16.168	2255	2258	2264	rVV2	38017	63363	0.79%	0.062%
67	16.268	2270	2275	2280	rVB	147433	226006	2.82%	0.222%
68	16.351	2286	2289	2292	rBV4	51048	71446	0.89%	0.070%
69	16.392	2293	2296	2299	rVB4	61917	78030	0.97%	0.077%
70	16.463	2306	2308	2315	rVB6	58521	84233	1.05%	0.083%
71	16.545	2319	2322	2328	rVV2	51815	96156	1.20%	0.094%
72	16.627	2332	2336	2342	rVV2	199102	355253	4.43%	0.349%
73	16.710	2346	2350	2356	rVV2	153560	314629	3.92%	0.309%
74	16.768	2356	2360	2362	rVV2	177021	261850	3.26%	0.257%
75	16.810	2362	2367	2371	rVV	2820565	4061080	50.64%	3.990%
76	16.851	2371	2374	2378	rVV	308608	474487	5.92%	0.466%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

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
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Title : SVOA CALIBRATION

77	16.910	2378	2384	2396	rVV2	3969521	6110885	76.20%	6.005%
78	17.004	2396	2400	2410	rVB5	168109	327512	4.08%	0.322%
79	17.204	2429	2434	2435	rBV	410971	556987	6.94%	0.547%
80	17.339	2454	2457	2466	rBV4	58627	124701	1.55%	0.123%
81	17.657	2509	2511	2515	rBV2	60707	70690	0.88%	0.069%
82	17.745	2518	2526	2534	rVV3	847900	1525458	19.02%	1.499%
83	17.815	2535	2538	2542	rVV3	121332	178153	2.22%	0.175%
84	17.880	2544	2549	2556	rVB	614675	931460	11.61%	0.915%
85	17.980	2563	2566	2569	rBV2	100789	148849	1.86%	0.146%
86	18.133	2589	2592	2599	rVB2	100371	181432	2.26%	0.178%
87	18.198	2600	2603	2606	rBV3	112966	151898	1.89%	0.149%
88	18.356	2627	2630	2633	rBV2	74012	86081	1.07%	0.085%
89	18.692	2684	2687	2692	rBV2	100601	129541	1.62%	0.127%
90	18.880	2715	2719	2724	rVB	1112663	1409117	17.57%	1.385%
91	19.045	2744	2747	2750	rBV	319286	312750	3.90%	0.307%
92	19.221	2769	2777	2781	rBV	5070913	7331116	91.41%	7.204%
93	19.639	2844	2848	2850	rBV	505178	706808	8.81%	0.695%
94	19.715	2858	2861	2866	rVB	128765	161854	2.02%	0.159%
95	20.333	2962	2966	2967	rBV	264029	336525	4.20%	0.331%
96	20.362	2967	2971	2980	rVB	760894	1250137	15.59%	1.228%
97	20.780	3039	3042	3048	rBV2	320995	362560	4.52%	0.356%
98	21.039	3081	3086	3092	rBV	3153624	4218323	52.60%	4.145%
99	23.568	3508	3516	3530	rVB	2928417	6446230	80.38%	6.334%
100	23.750	3539	3547	3555	rBV	1857886	4209543	52.49%	4.136%

Sum of corrected areas: 101770927

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
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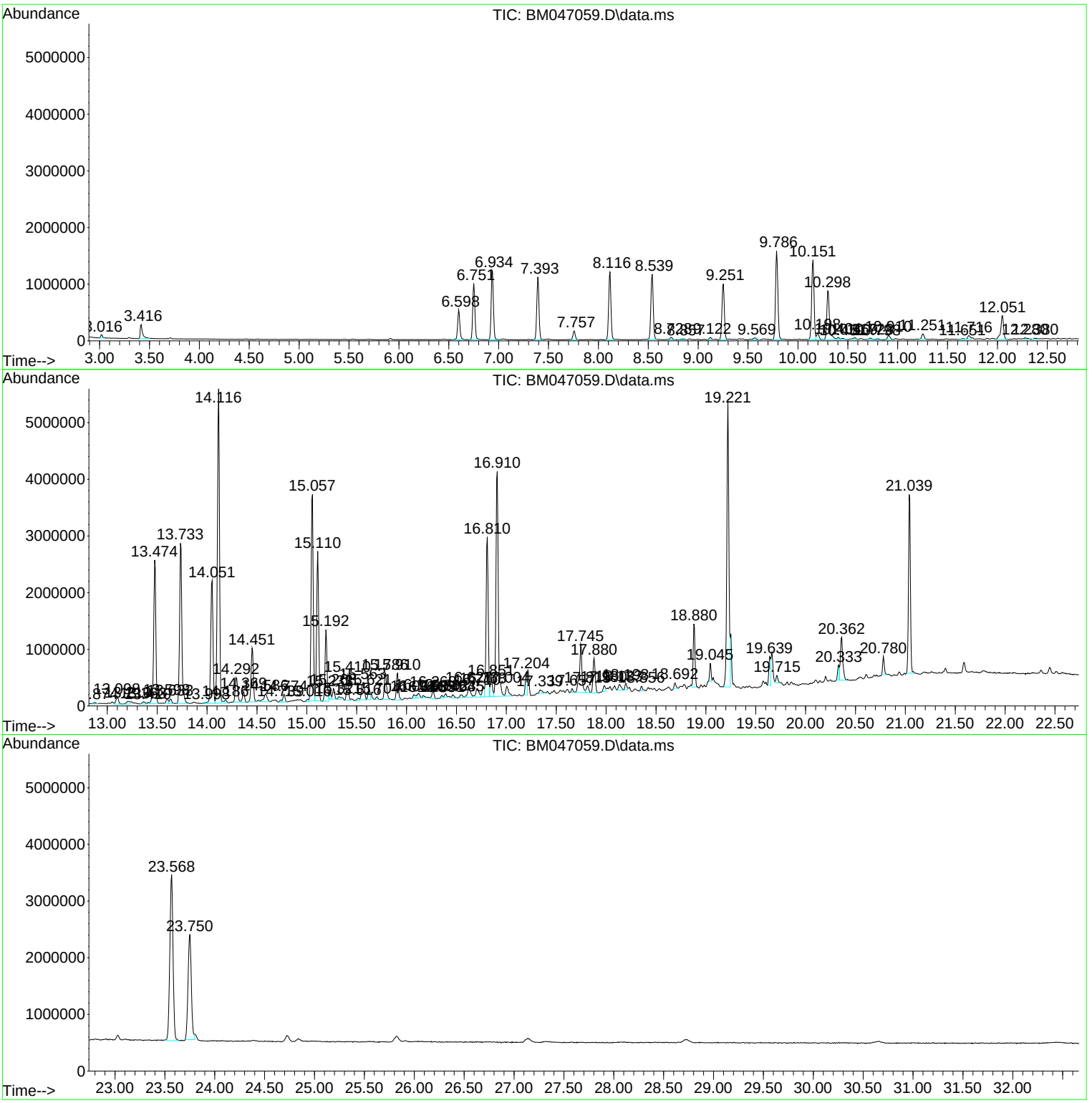
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0B10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

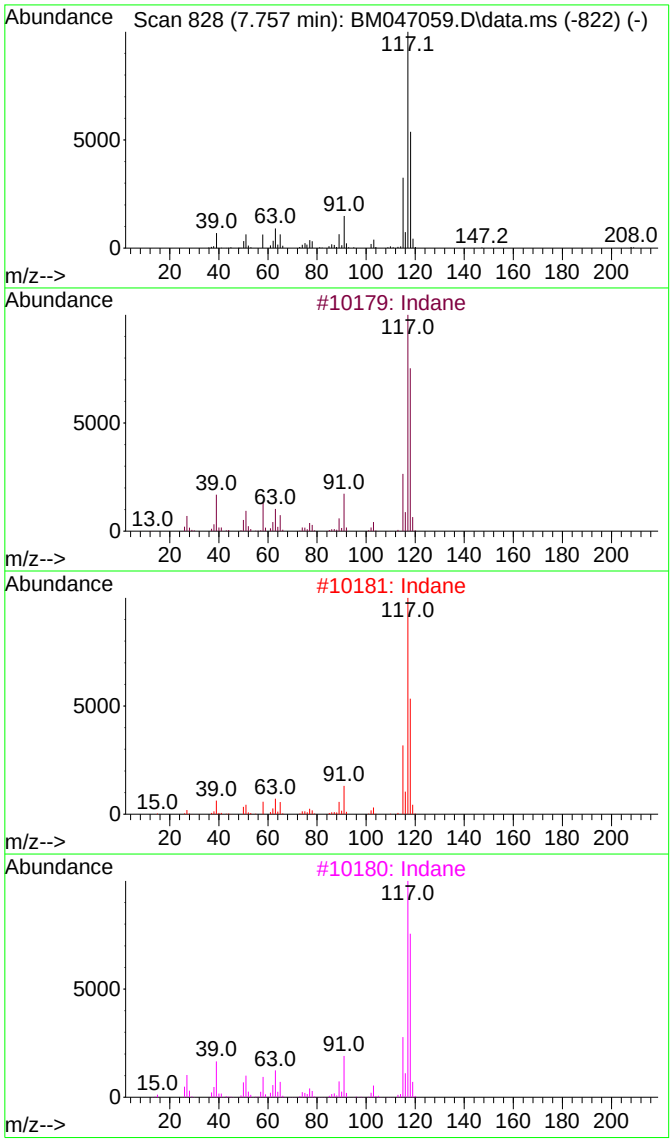
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

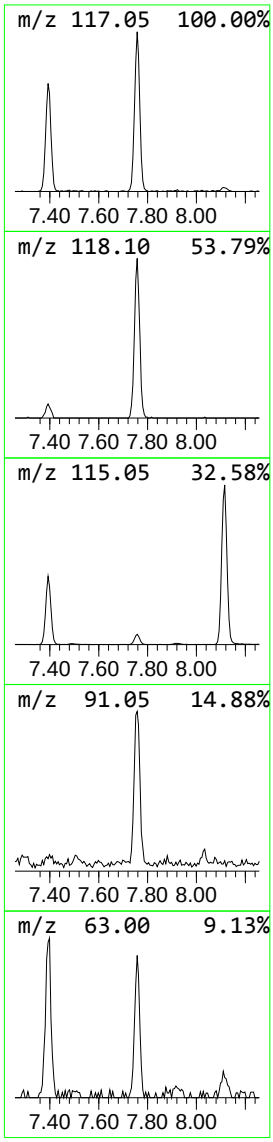
****

Peak Number 1 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.757	2.92 ng/ul	244515	1,4-Dichlorobenzene-d4	7.393

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	90
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	87
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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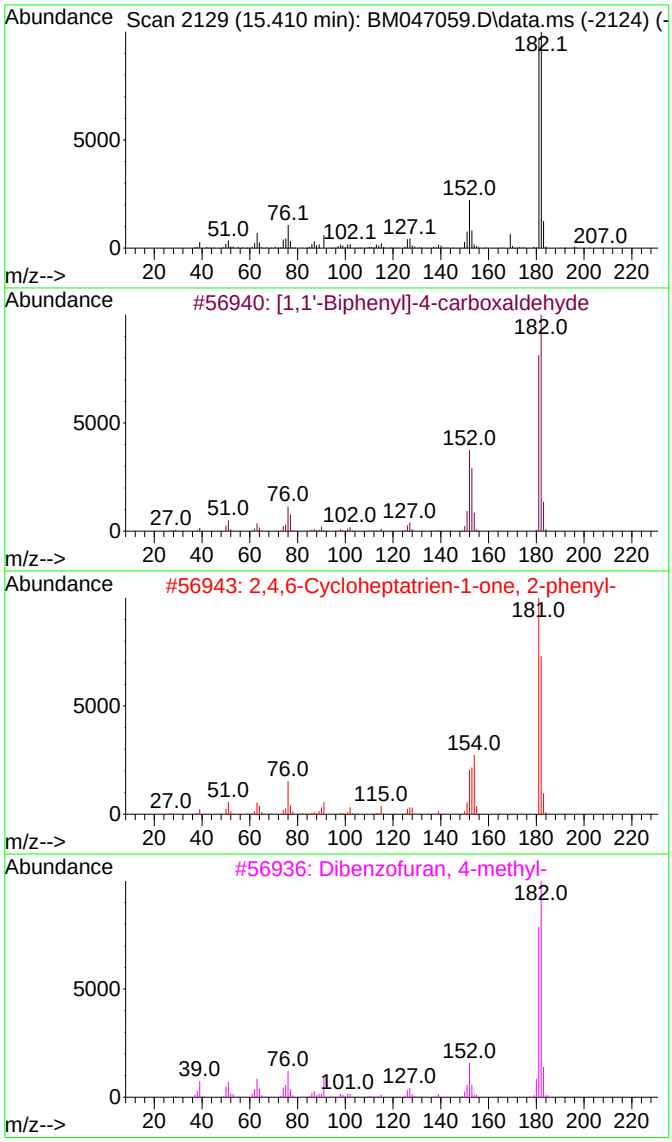
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

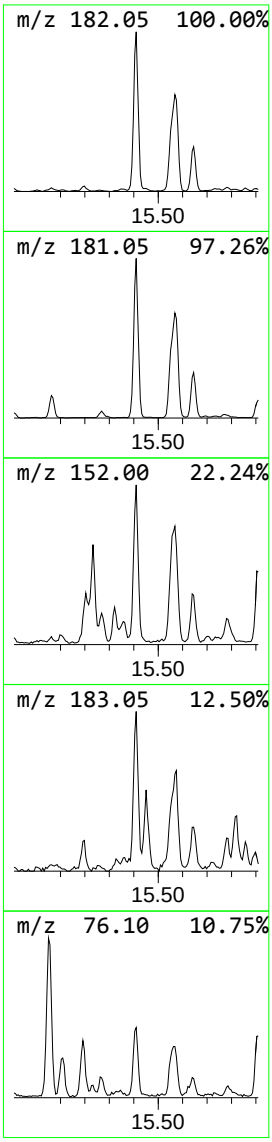
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.410	3.54 ng/ul	592288	Acenaphthene-d10	14.051

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
2	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
3	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	86
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
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Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

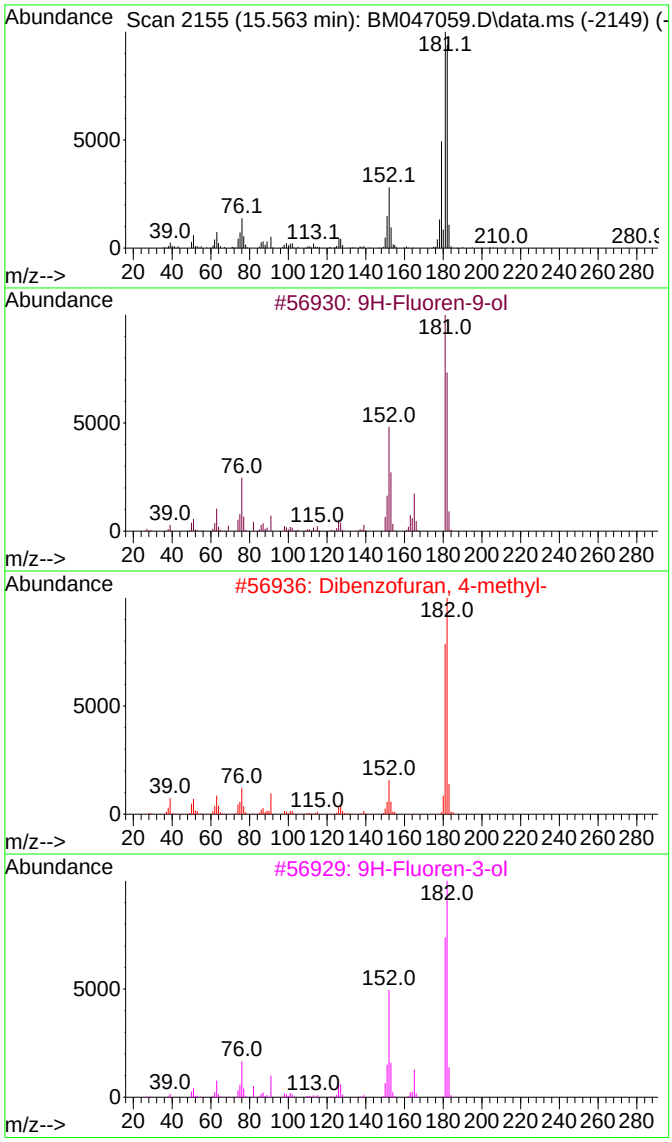
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

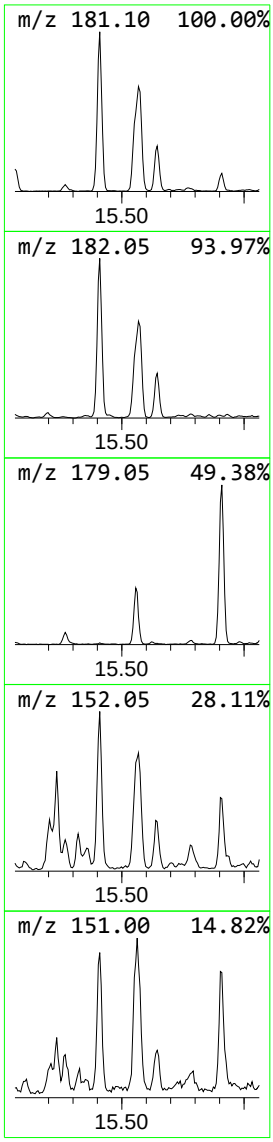
****

Peak Number 3 9H-Fluoren-9-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.563	3.12 ng/ul	633409	Phenanthrene-d10	16.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
2	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	64
3	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	64
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	62
5	2-Hydroxyfluorene	182	C13H10O	002443-58-5	60





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Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

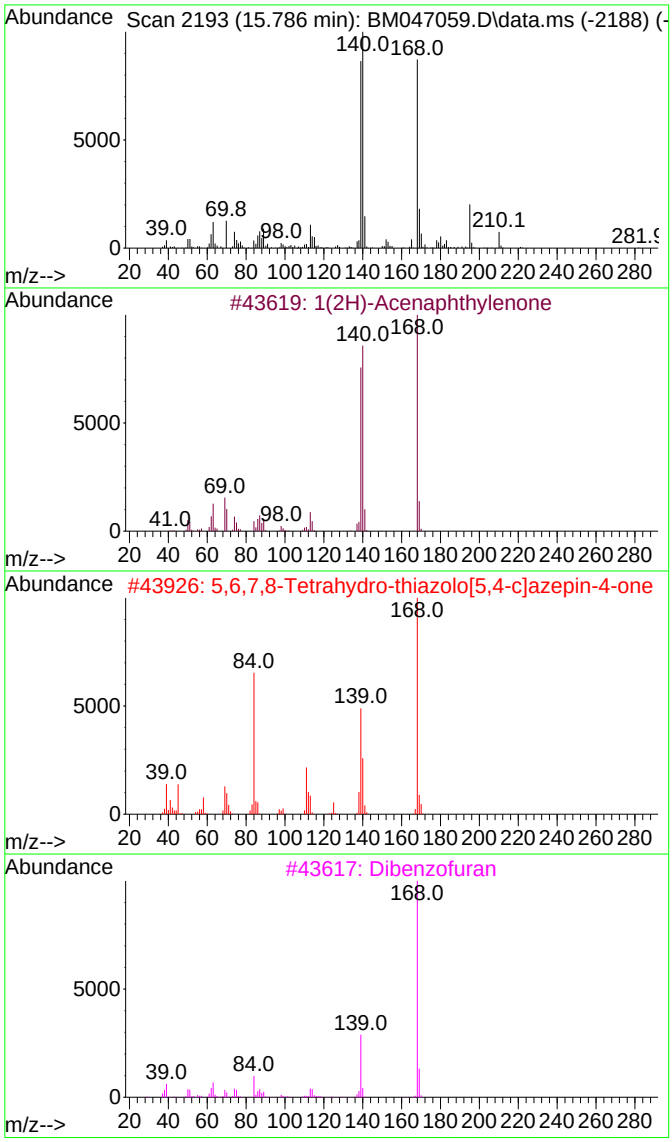
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

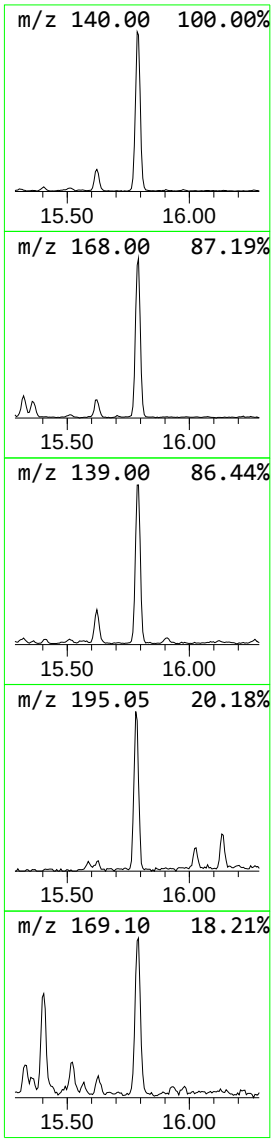
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1(2H)-Acenaphthylenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.786	3.80 ng/ul	772238	Phenanthrene-d10	16.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	93
2	5,6,7,8-Tetrahydro-thiazolo[5,4-...	168	C7H8N2OS	1000317-75-4	50
3	Dibenzofuran	168	C12H8O	000132-64-9	49
4	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	49
5	7(4H)-Benzothiazolone, 2-amino-5...	168	C7H8N2OS	017583-10-7	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

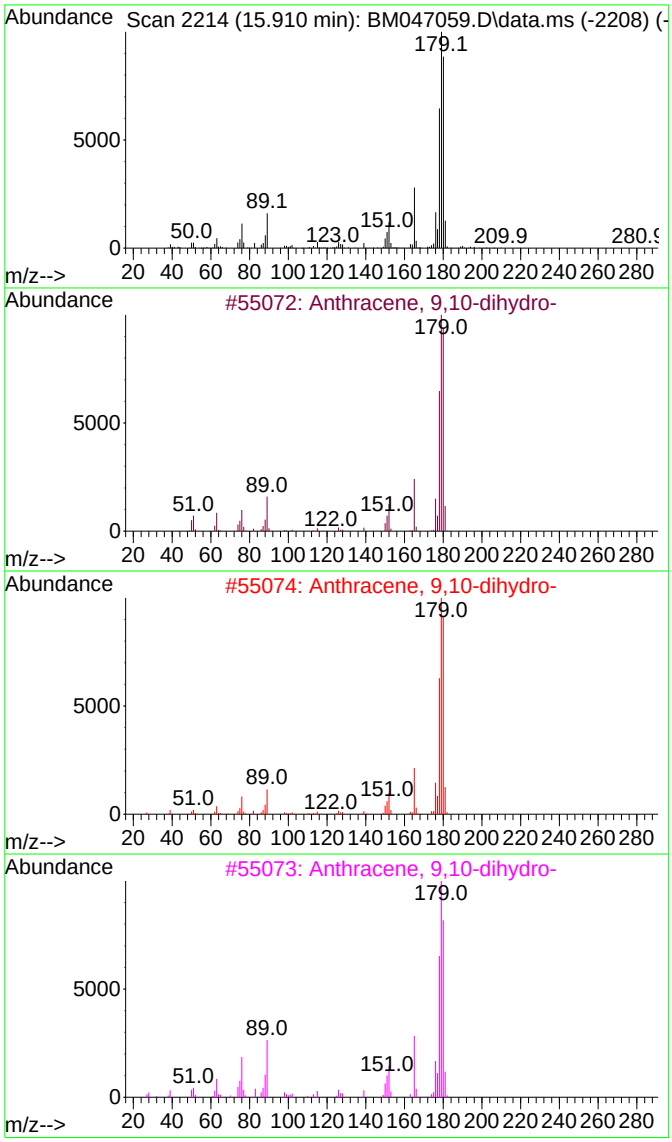
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

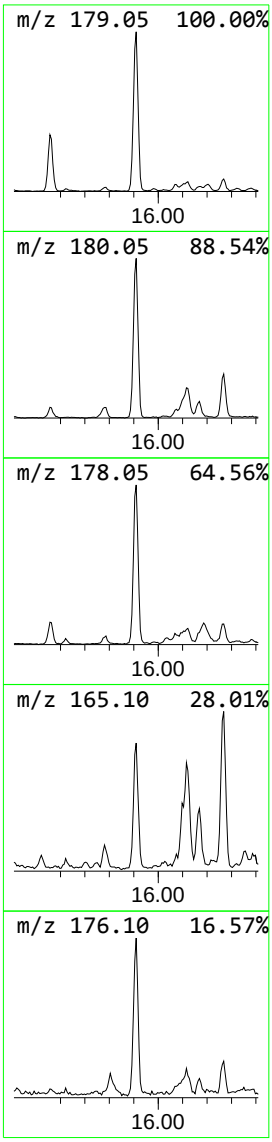
****

Peak Number 5 Anthracene, 9,10-dihydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.910	3.39 ng/ul	688376	Phenanthrene-d10	16.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	97
2	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	97
3	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	96
4	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	96
5	Phenanthrene, 9,10-dihydro-	180	C14H12	000776-35-2	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

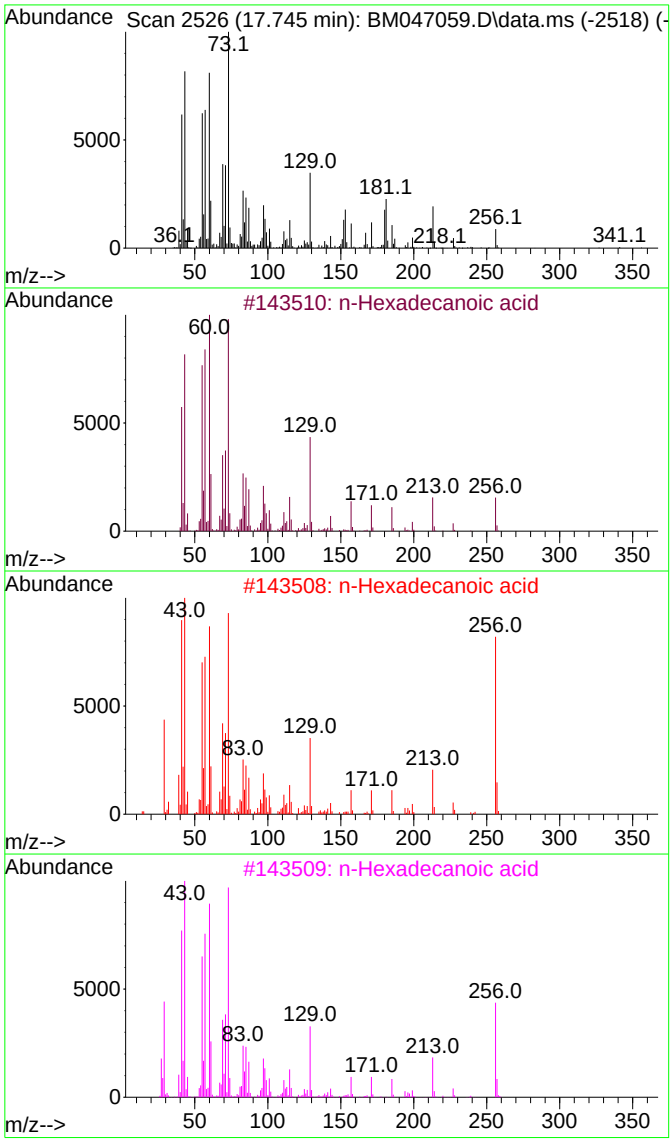
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

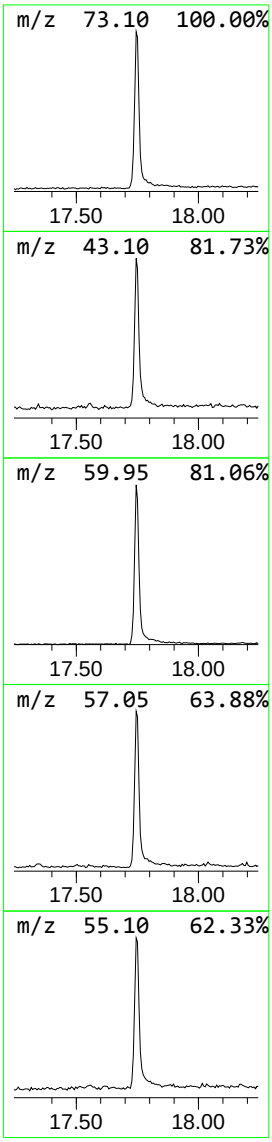
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.745	7.51 ng/ul	1525460	Phenanthrene-d10	16.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

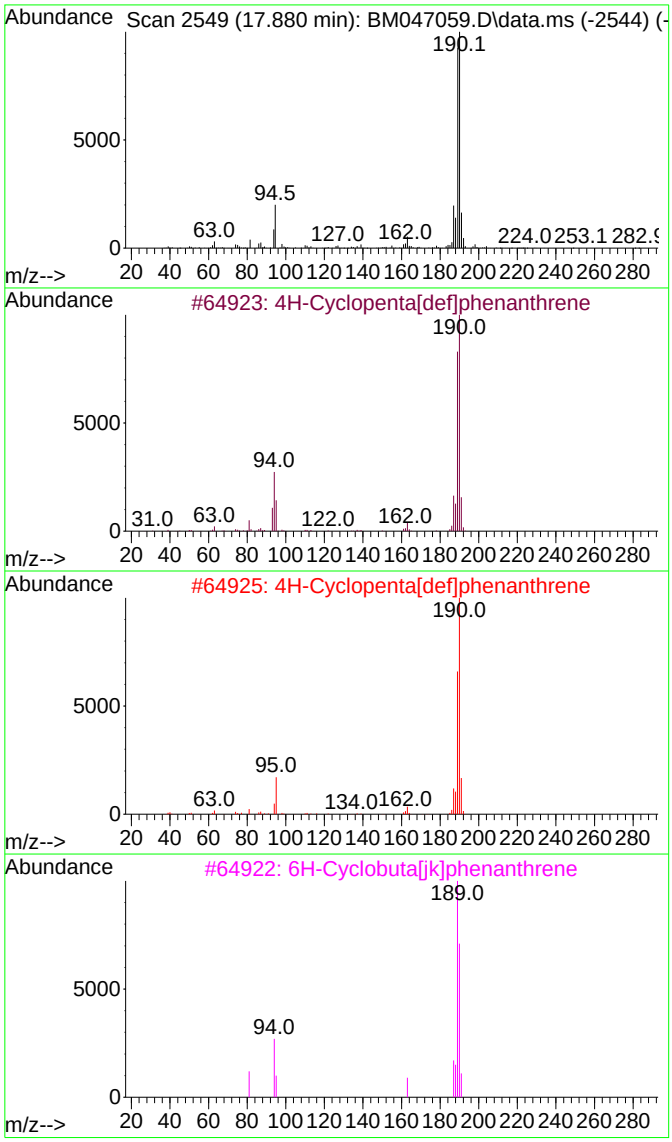
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

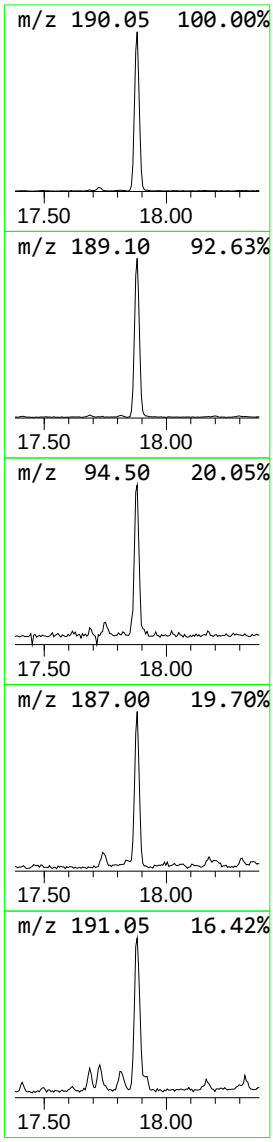
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.880	4.59 ng/ul	931460	Phenanthrene-d10	16.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	97
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	64
5	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

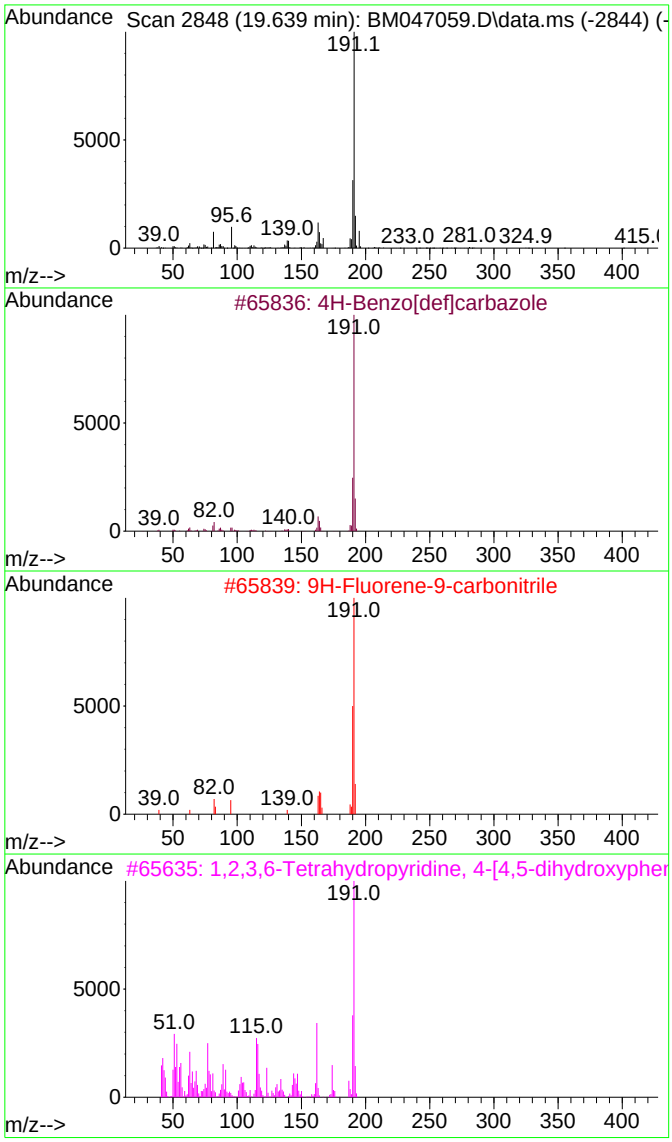
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

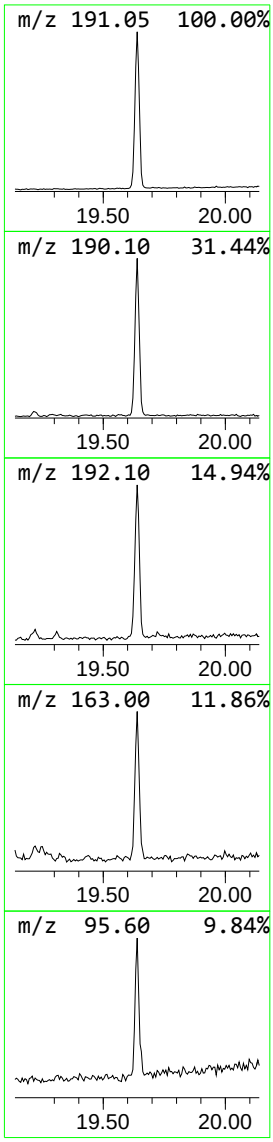
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4H-Benzo[def]carbazole Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.639	3.35 ng/ul	706808	Chrysene-d12	21.039

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Benzo[def]carbazole	191	C14H9N	000203-65-6	94
2	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	59
3	1,2,3,6-Tetrahydropyridine, 4-[4...	191	C11H13NO2	090684-17-6	59
4	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	58
5	9H-Fluorene-2-carbonitrile	191	C14H9N	002523-48-0	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

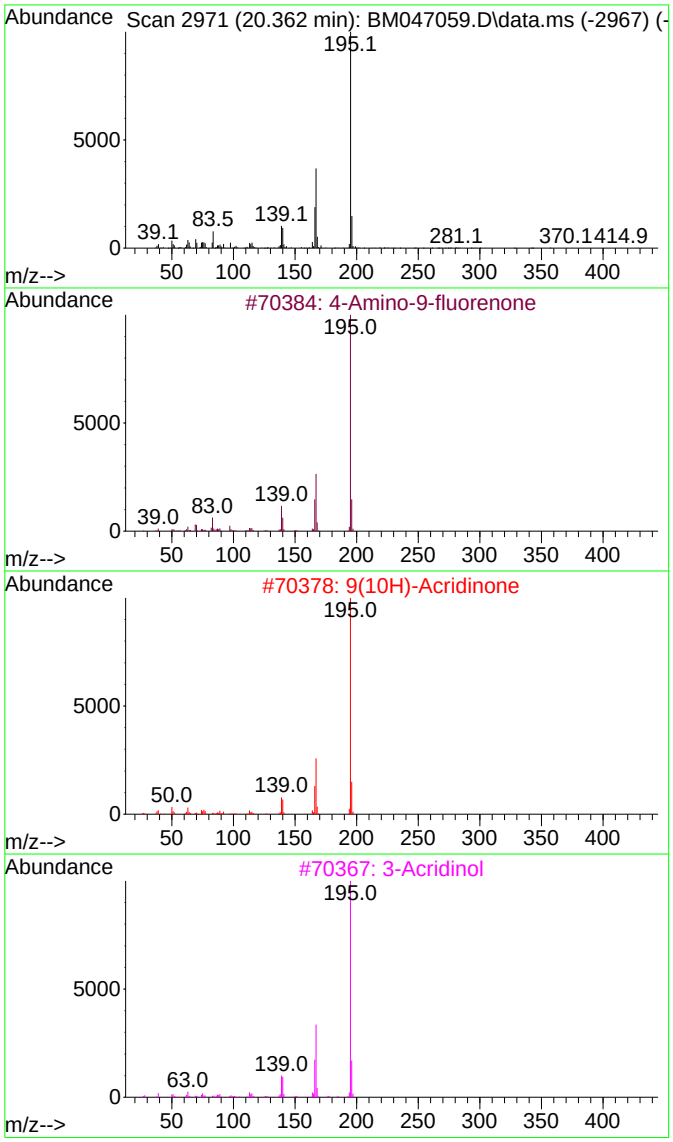
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

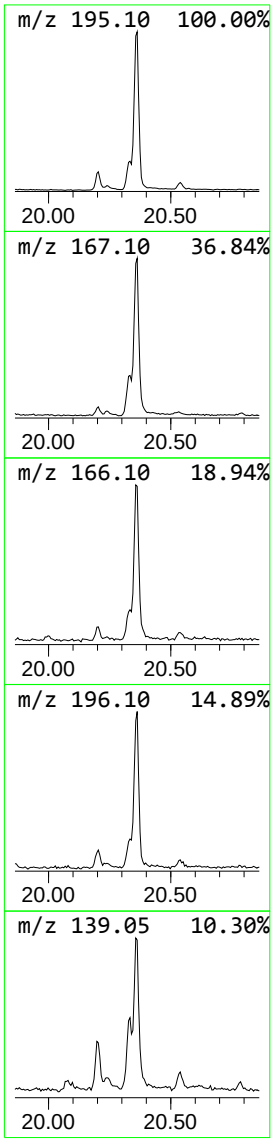
****

Peak Number 9 4-Amino-9-fluorenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.362	5.93 ng/ul	1250140	Chrysene-d12	21.039

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	95
2	9(10H)-Acridinone	195	C13H9NO	000578-95-0	94
3	3-Acridinol	195	C13H9NO	007132-70-9	93
4	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	90
5	9(10H)-Acridinone	195	C13H9NO	000578-95-0	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080624\
Data File : BM047059.D
Acq On : 07 Aug 2024 16:02
Operator : RC/JU
Sample : P3378-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.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
Indane	7.757	2.9	ng/ul	244515	1	7.393	1672790	20.0
[1,1'-Biphenyl]...	15.410	3.5	ng/ul	592288	3	14.051	3346020	20.0
9H-Fluoren-9-ol	15.563	3.1	ng/ul	633409	4	16.810	4061080	20.0
1(2H)-Acenaphth...	15.786	3.8	ng/ul	772238	4	16.810	4061080	20.0
Anthracene, 9,1...	15.910	3.4	ng/ul	688376	4	16.810	4061080	20.0
n-Hexadecanoic ...	17.745	7.5	ng/ul	1525460	4	16.810	4061080	20.0
4H-Cyclopenta[d...	17.880	4.6	ng/ul	931460	4	16.810	4061080	20.0
4H-Benzo[def]ca...	19.639	3.4	ng/ul	706808	5	21.039	4218320	20.0
4-Amino-9-fluor...	20.362	5.9	ng/ul	1250140	5	21.039	4218320	20.0