LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : E05Y6DL

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: BP021474.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.634	782	790	817	rBV2	208486	845454	12.89%	1.678%
2	10.375	1246	1256	1260	rBV	540947	1056525	16.11%	2.097%
3	10.428	1260	1265	1286	rVB	588130	1608855	24.53%	3.193%
4	11.963	1521	1526	1531	rBV3	6634	10772	0.16%	0.021%
5	12.069	1537	1544	1566	rBV	373013	949496	14.48%	1.884%
6	12.281	1573	1580	1595	rVB	159368	425357	6.49%	0.844%
7	13.093	1712	1718	1731	rBV	112899	241167	3.68%	0.479%
8	13.281	1738	1750	1765	rVB	50896	119999	1.83%	0.238%
9	13.440	1767	1777	1794	rBV3	59368	201677	3.08%	0.400%
10	13.581	1794	1801	1806	rBV	90673	181445	2.77%	0.360%
11	13.628	1806	1809	1819	rVB	56449	94276	1.44%	0.187%
12	13.828	1834	1843	1855	rBV	38416	99948	1.52%	0.198%
13	13.981	1862	1869	1878	rVB4	25720	52538	0.80%	0.104%
14	14.251	1908	1915	1922	rBV	1086087	1871760	28.54%	3.715%
15	14.322	1922	1927	1934	rVB	870471	1509454	23.02%	2.996%
16	14.492	1948	1956	1964	rBV4	20622	62142	0.95%	0.123%
17	14.581	1964	1971	1977	rVB2	30271	53208	0.81%	0.106%
18	14.675	1977	1987	1999	rBV	544895	1049429	16.00%	2.083%
19	14.763	1999	2002	2007	rVV	28881	43003	0.66%	0.085%
20	14.834	2010	2014	2018	rVV5	4982	8936	0.14%	0.018%
21	14.887	2018	2023	2030	rVB4	14191	27256	0.42%	0.054%
22	14.957	2030	2035	2044	rVB4	13281	27061	0.41%	0.054%
23	15.075	2047	2055	2058	rBV5	10911	23089	0.35%	0.046%
24	15.110	2058	2061	2067	rVB4	10528	13843	0.21%	0.027%
25	15.228	2075	2081	2084	rBV6	6891	11533	0.18%	0.023%
26	15.275	2084	2089	2092	rVV4	17756	27839	0.42%	0.055%
27	15.334	2092	2099	2113	rVV	777377	1545123	23.56%	3.066%
28	15.463	2114	2121	2124	rVV4	47838	132040	2.01%	0.262%
29	15.498	2124	2127	2133	rVV2	97429	166223	2.53%	0.330%
30	15.545	2133	2135	2139	rVV3	19982	32879	0.50%	0.065%
31	15.592	2139	2143	2148	rVV	52520	93516	1.43%	0.186%
32	15.651	2148	2153	2161	rVB2	83726	175018	2.67%	0.347%
33	15.804	2174	2179	2191	rBV	105866	252229	3.85%	0.501%
34	15.892	2191	2194	2206	rVV5	21607	54278	0.83%	0.108%
35	15.998	2206	2212	2224	rVB7	20957	50139	0.76%	0.100%
36	16.157	2235	2239	2245	rBV2	45609	62931	0.96%	0.125%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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	16.369	2263	2275	2281	rBV	81068	175485	2.68%	0.348%
38	16.428	2281	2285	2289	rVB2	23469	38634	0.59%	0.077%
39	16.522	2297	2301	2306	rVB2	32826	44206	0.67%	0.088%
40	16.569	2306	2309	2311	rBV3	14937	19248	0.29%	0.038%
41	16.598	2311	2314	2317	rVV3	28798	40254	0.61%	0.080%
42	16.634	2317	2320	2324	rVV4	23928	38935	0.59%	0.077%
43	16.798	2343	2348	2349	rBV2	31857	41052	0.63%	0.081%
44	16.898	2360	2365	2379	rVB	274531	411027	6.27%	0.816%
45	17.069	2387	2394	2397	rBV	1520811	2174857	33.17%	4.316%
46	17.116	2397	2402	2411	rVV	3629877	6557641	100.00%	13.014%
47	17.222	2414	2420	2432	rVV	1050915	2300586	35.08%	4.566%
48	17.322	2433	2437	2445	rVV2	71513	161581	2.46%	0.321%
49	17.392	2445	2449	2455	rVB	27517	45338	0.69%	0.090%
50	17.533	2467	2473	2483	rBV	355208	698003	10.64%	1.385%
51	17.616	2483	2487	2491	rVV	44020	66668	1.02%	0.132%
52	17.657	2491	2494	2496	rVV4	9639	14543	0.22%	0.029%
53	17.692	2496	2500	2508	rVB5	29529	62891	0.96%	0.125%
54	17.992	2545	2551	2555	rBV	188059	322570	4.92%	0.640%
55	18.051	2555	2561	2569	rVV	224162	449666	6.86%	0.892%
56	18.139	2570	2576	2579	rVV	90026	170679	2.60%	0.339%
57	18.186	2579	2584	2589	rVV	471367	798670	12.18%	1.585%
58	18.233	2589	2592	2598	rVB3	98759	165317	2.52%	0.328%
59	18.333	2606	2609	2614	rVB	14528	20047	0.31%	0.040%
60	18.386	2614	2618	2624	rBV3	20081	39970	0.61%	0.079%
61	18.510	2635	2639	2649	rVB	175563	237662	3.62%	0.472%
62	18.628	2657	2659	2666	rVB5	19079	27388	0.42%	0.054%
63	18.763	2678	2682	2687	rBV4	38883	52605	0.80%	0.104%
64	18.816	2688	2691	2695	rBV	30770	36261	0.55%	0.072%
65	18.863	2696	2699	2703	rBV3	19773	30169	0.46%	0.060%
66	18.939	2708	2712	2715	rBV	53919	76828	1.17%	0.152%
67	19.004	2721	2723	2726	rVB2	24848	23082	0.35%	0.046%
68	19.216	2753	2759	2768	rBV	3242455	4667960	71.18%	9.264%
69	19.475	2798	2803	2811	rBV2	82550	147592	2.25%	0.293%
70	19.592	2812	2823	2835	rBV	2177493	4039237	61.60%	8.016%
71	19.686	2835	2839	2845	rVB	85116	140833	2.15%	0.279%
72	19.798	2854	2858	2862	rBV	128605	177052	2.70%	0.351%
73	19.898	2871	2875	2877	rBV	30272	49087	0.75%	0.097%
74	19.939	2878	2882	2884	rBV2	69917	110073	1.68%	0.218%
75	20.127	2909	2914	2921	rVB	318382	516096	7.87%	1.024%
76	20.222	2926	2930	2935	rVV	424842	576483	8.79%	1.144%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
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Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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

77	20.280	2936	2940	2951	rVB4	128087	329567	5.03%	0.654%
78	20.439	2963	2967	2970	rBV	54003	88305	1.35%	0.175%
79	21.098	3075	3079	3081	rBV	90654	131292	2.00%	0.261%
80	21.127	3082	3084	3090	rVB	101770	140019	2.14%	0.278%
81	21.186	3091	3094	3097	rBV	79394	112481	1.72%	0.223%
82	21.510	3141	3149	3155	rBV3	1494748	3879883	59.17%	7.700%
83	21.569	3155	3159	3173	rVB	645255	1456562	22.21%	2.891%
84	21.716	3181	3184	3190	rVB	166039	289908	4.42%	0.575%
85	23.774	3526	3534	3540	rBV	463158	1395583	21.28%	2.770%
86	24.492	3651	3656	3663	rVB	144434	306845	4.68%	0.609%
87	24.645	3676	3682	3692	rVB2	304766	742874	11.33%	1.474%
88	24.792	3699	3707	3719	rBV	926485	2569191	39.18%	5.099%

Sum of corrected areas: 50389224

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
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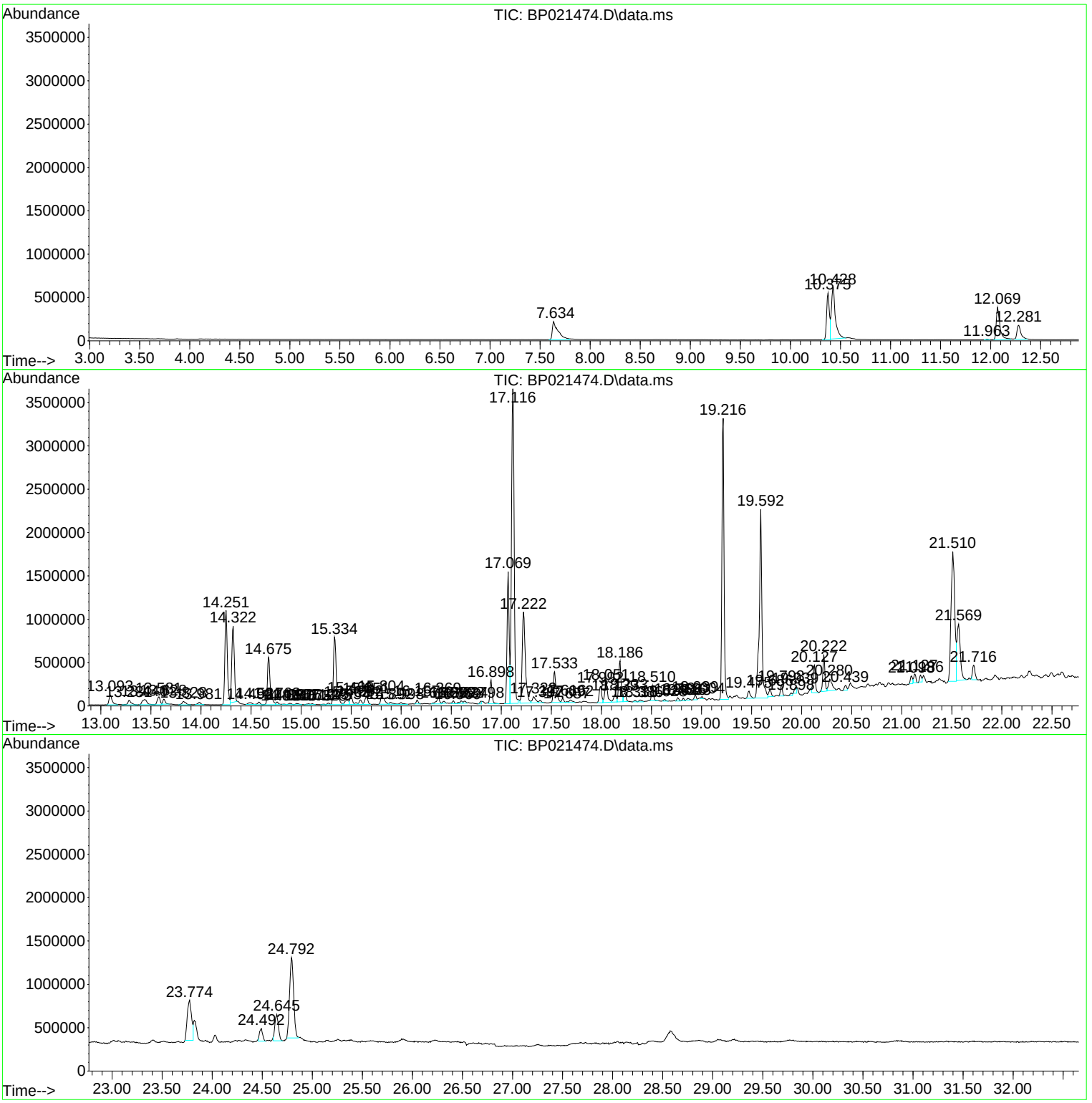
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

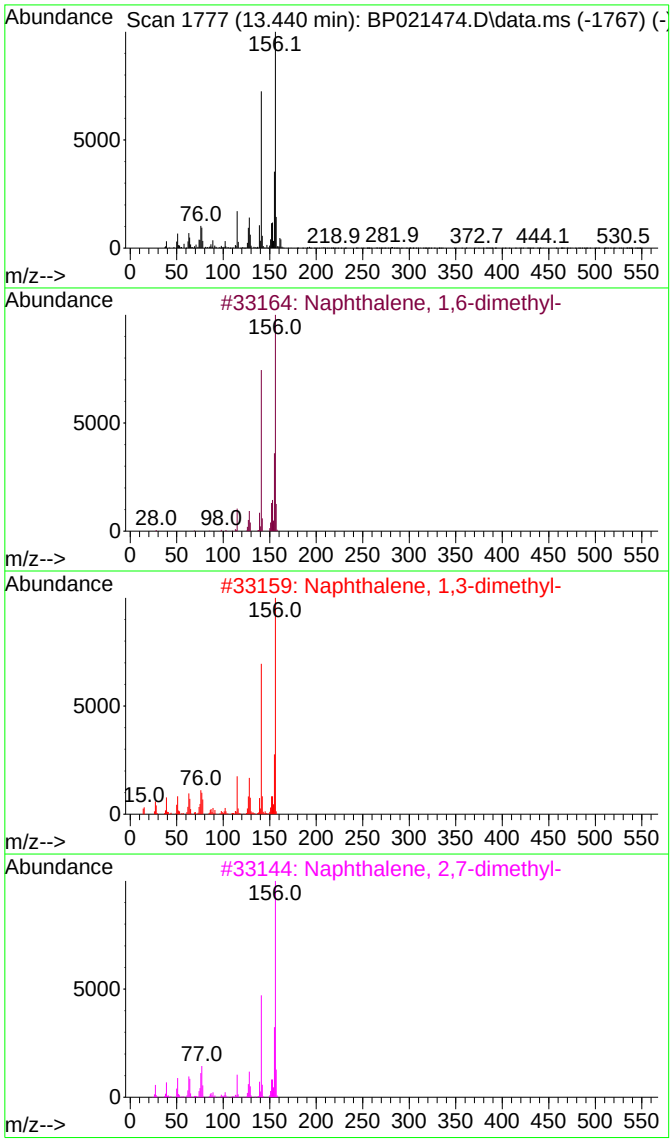
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

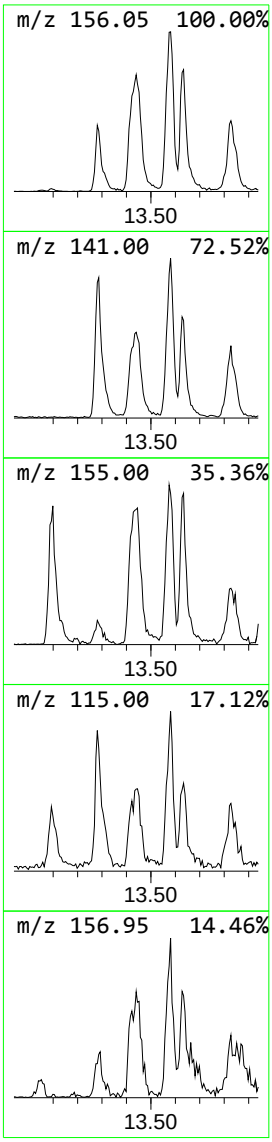
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Naphthalene, 1,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.440	2.15 ng/ul	201677	Acenaphthene-d10	14.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

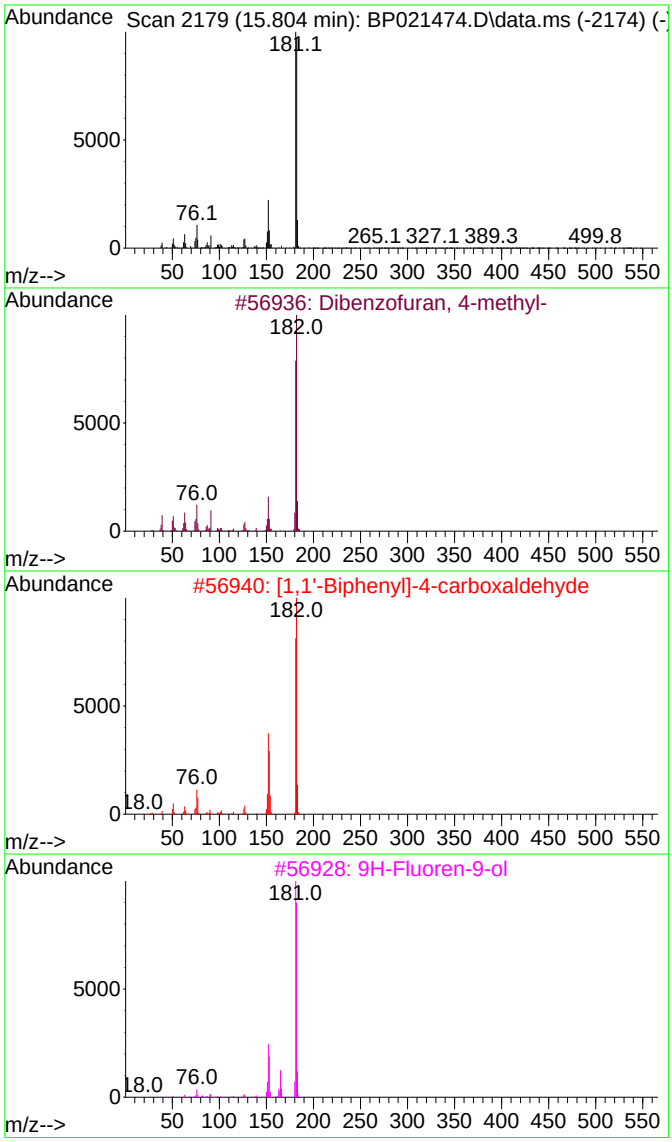
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

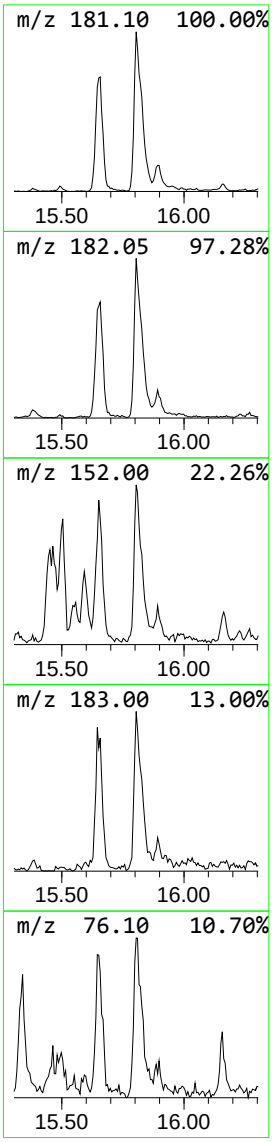
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Dibenzofuran, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	2.32 ng/ul	252229	Phenanthrene-d10	17.069

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	93
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	91
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	90
4	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	87
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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

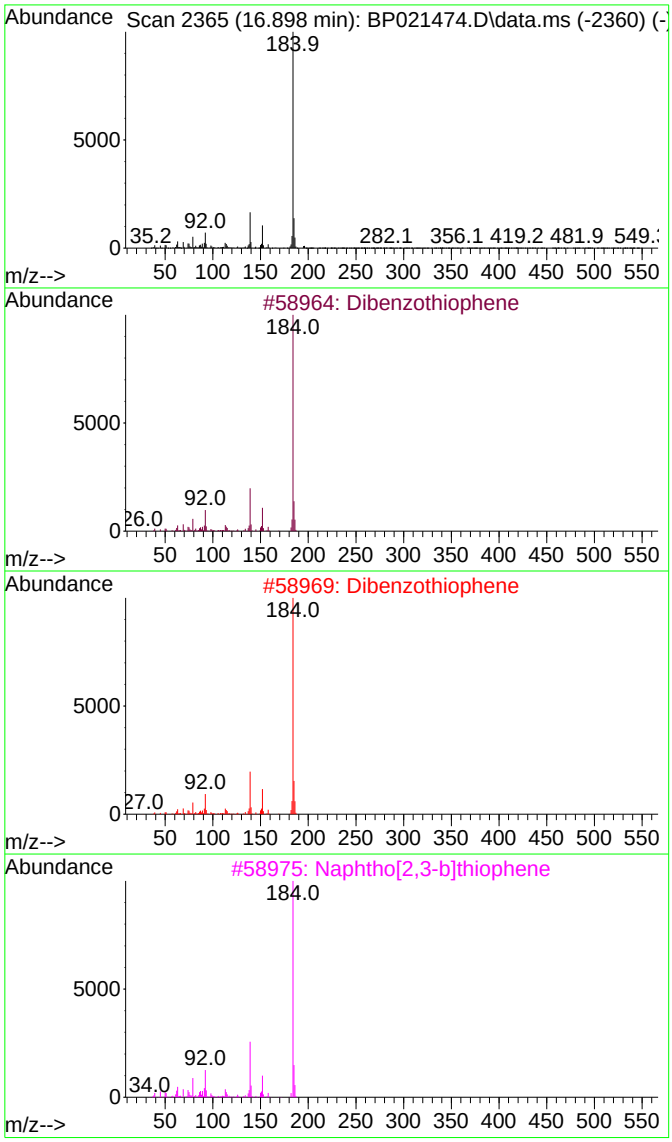
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

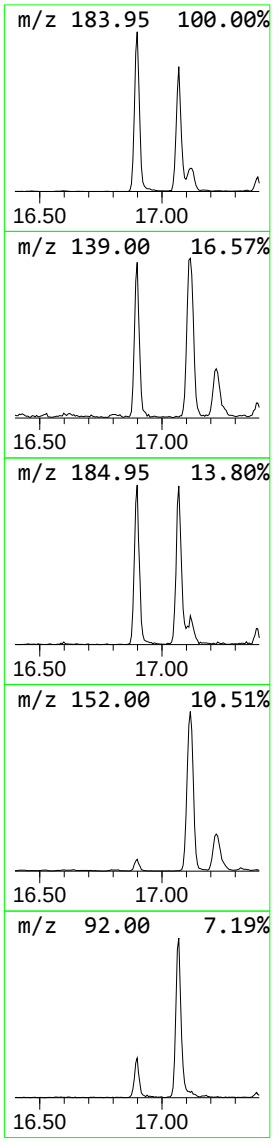
****

Peak Number 3 Dibenzothiophene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.898	3.78 ng/ul	411027	Phenanthrene-d10	17.069

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	97
2	Dibenzothiophene	184	C12H8S	000132-65-0	97
3	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	97
4	Dibenzothiophene	184	C12H8S	000132-65-0	96
5	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

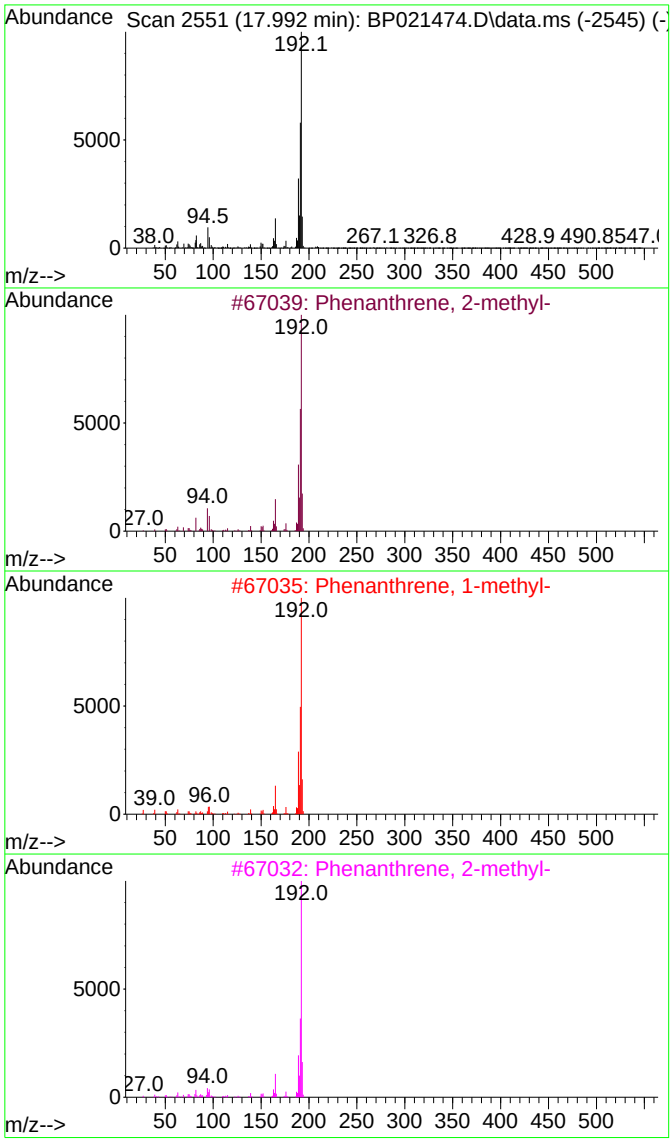
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

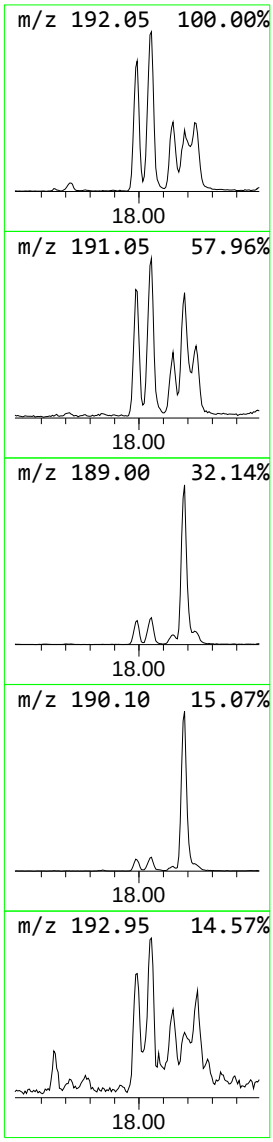
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	2.97 ng/ul	322570	Phenanthrene-d10	17.069

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

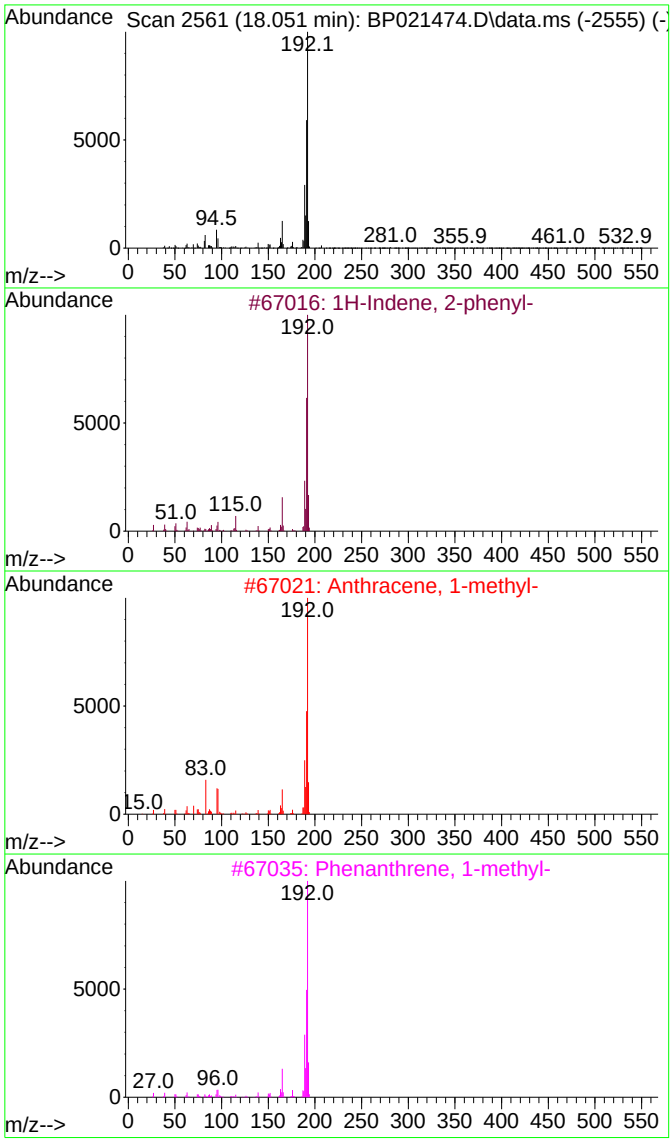
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

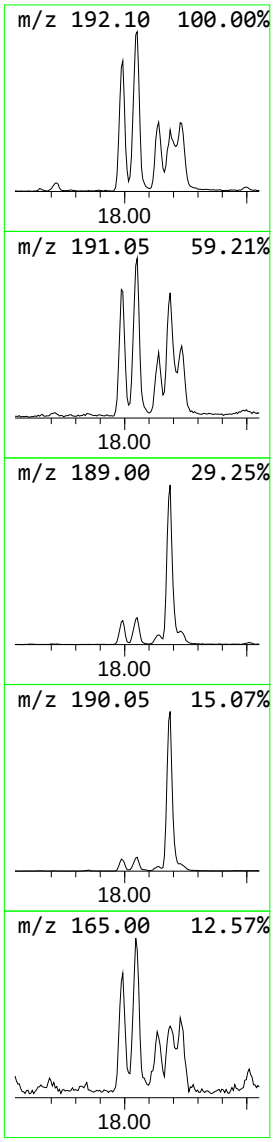
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1H-Indene, 2-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	4.14 ng/ul	449666	Phenanthrene-d10	17.069

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

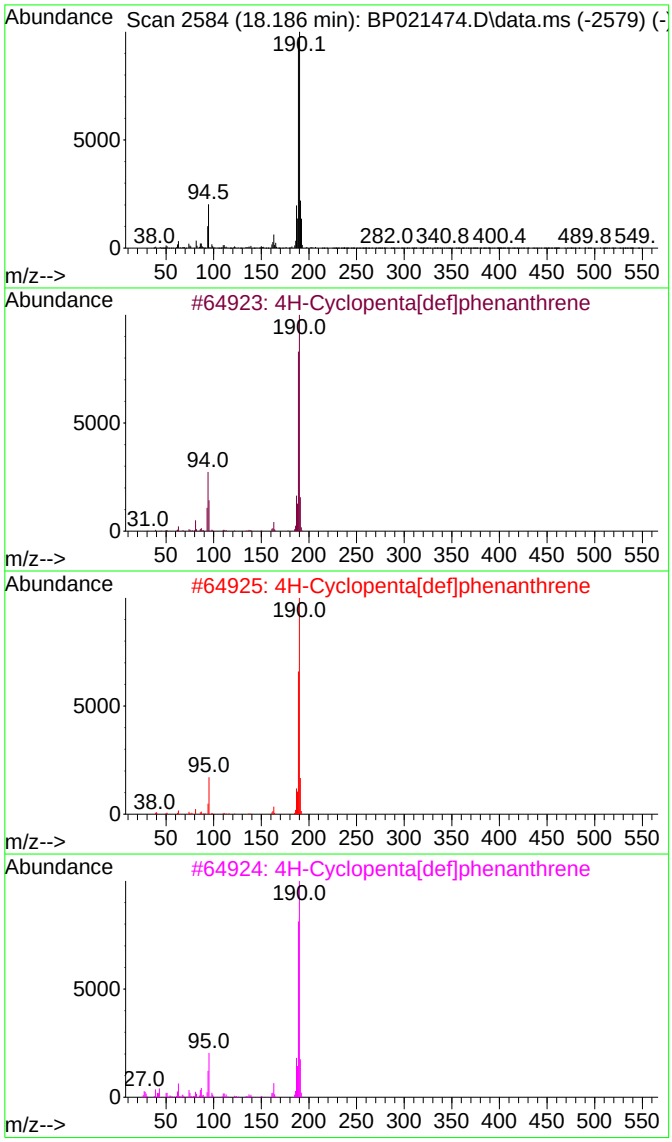
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

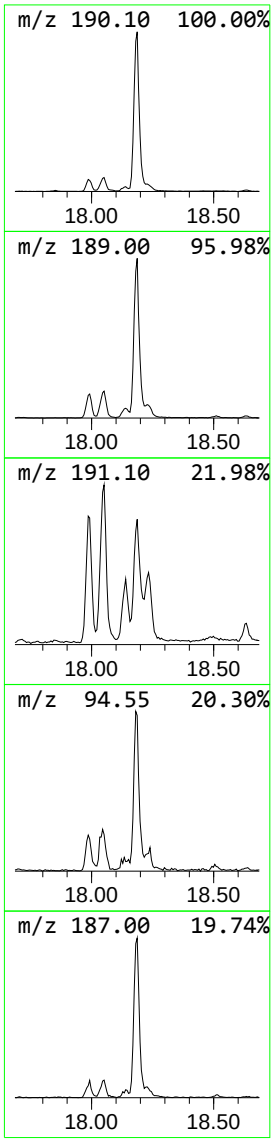
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	7.34 ng/ul	798670	Phenanthrene-d10	17.069

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	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
4	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	59
5	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	56





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Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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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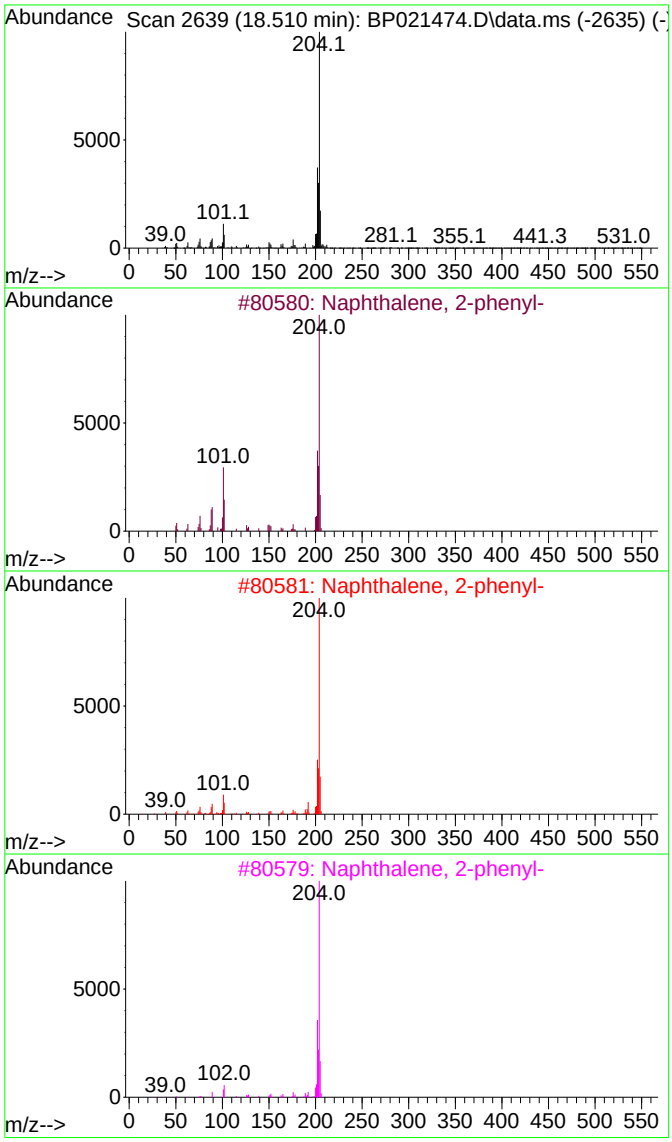
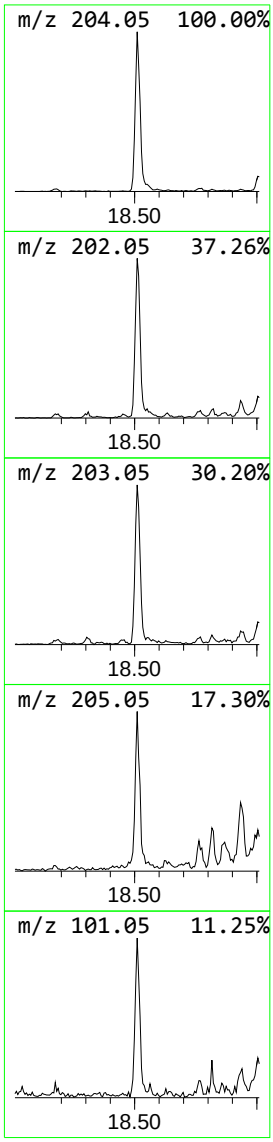
Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.510	2.19 ng/ul	237662	Phenanthrene-d10	17.069

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
4	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	90
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89





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Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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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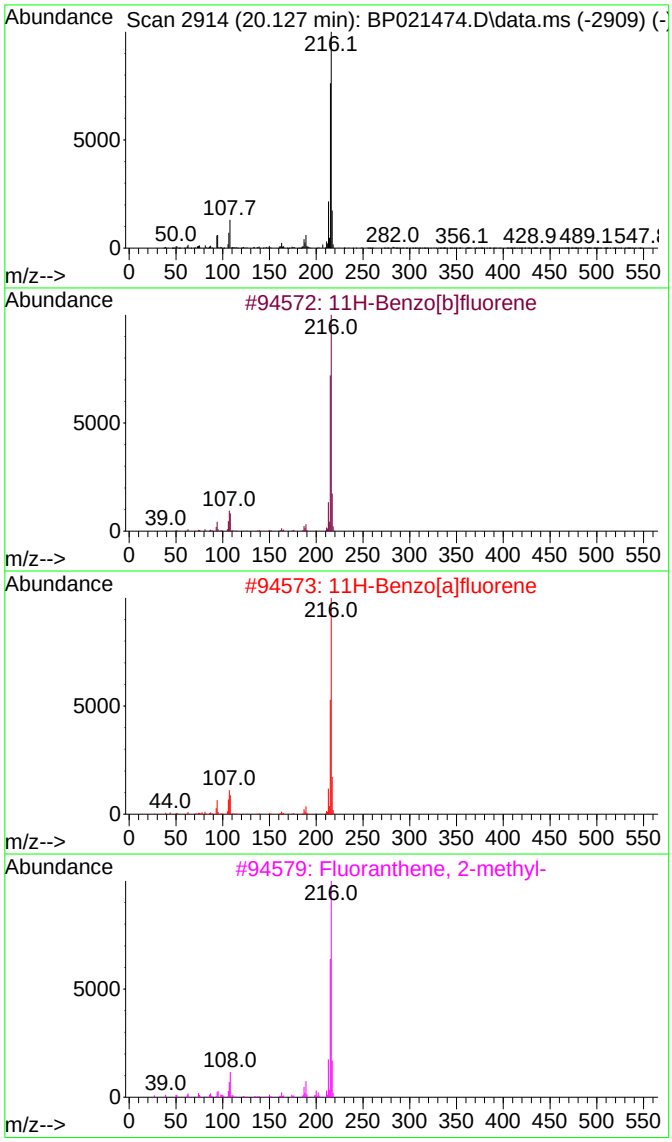
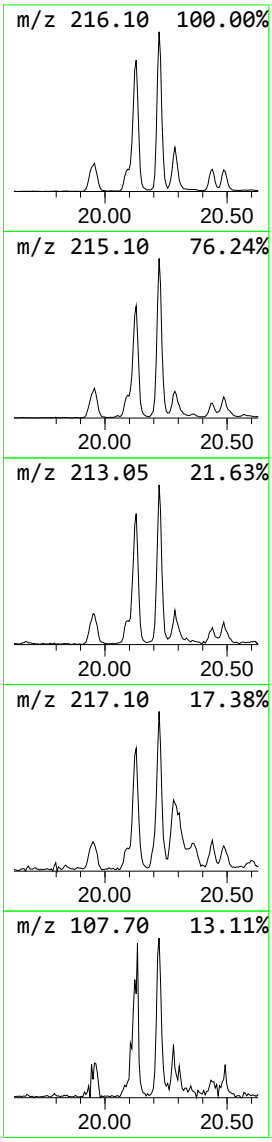
Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.127	2.66 ng/ul	516096	Chrysene-d12	21.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91





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Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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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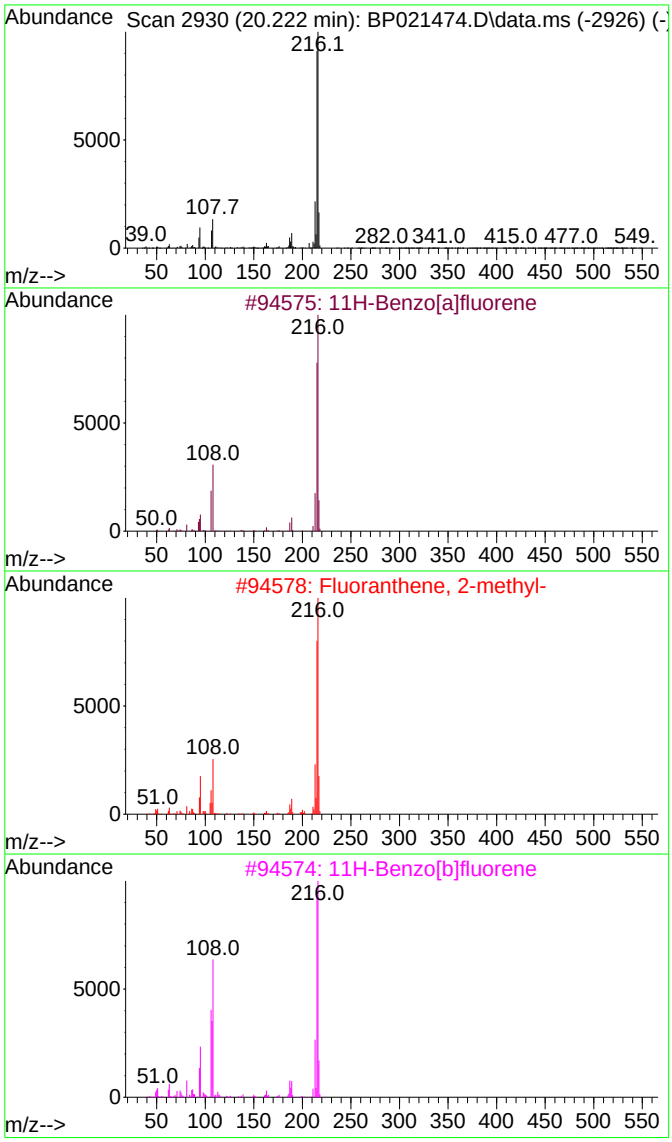
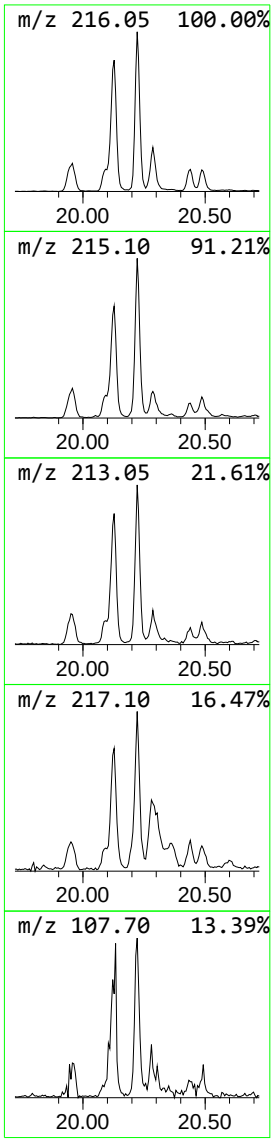
Peak Number 9 11H-Benzo[a]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.221	2.97 ng/ul	576483	Chrysene-d12	21.516

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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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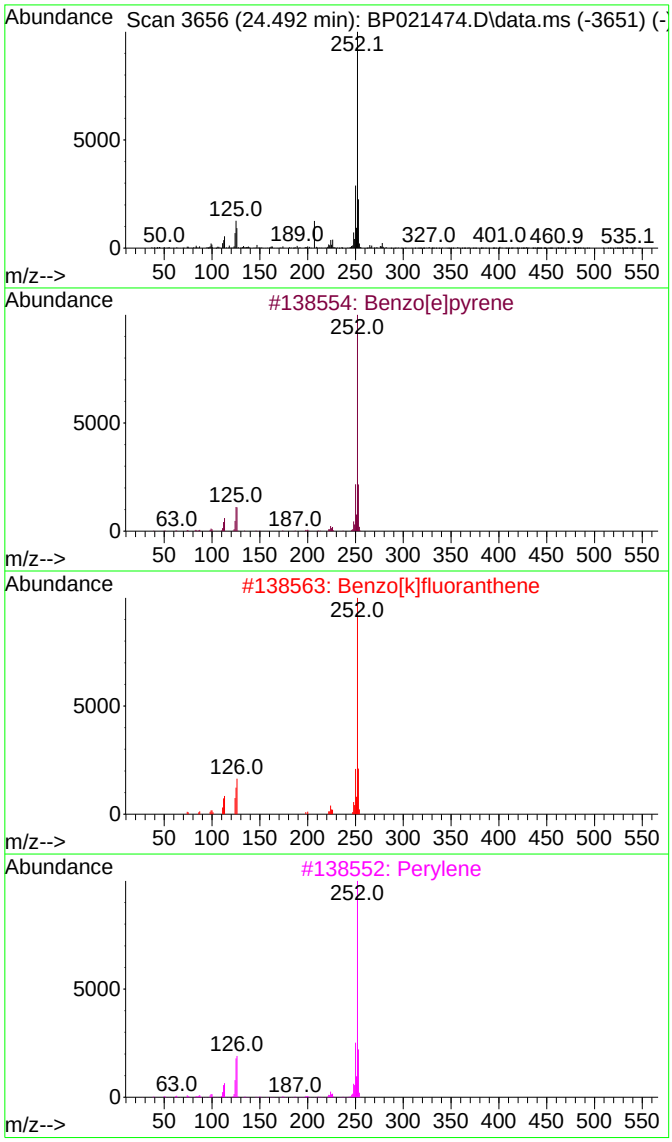
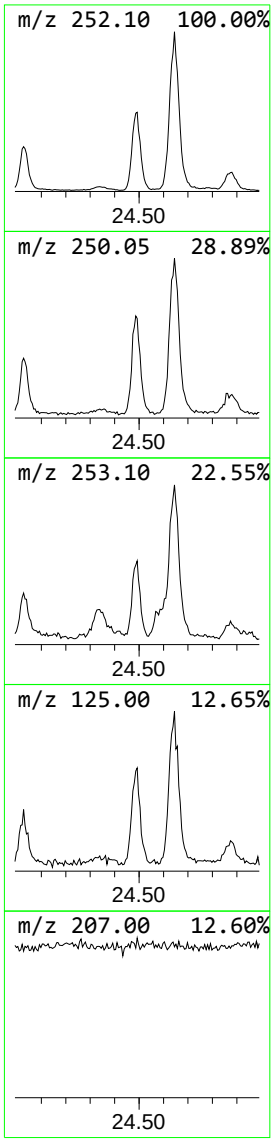
Peak Number 10 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.492	2.39 ng/ul	306845	Perylene-d12	24.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3	Perylene	252	C20H12	000198-55-0	95
4	Perylene	252	C20H12	000198-55-0	95
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
Data File : BP021474.D
Acq On : 09 Aug 2024 17:17
Operator : RC/JU
Sample : P3435-14DL 100X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E05Y6DL

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Naphthalene, 1,...	13.440	2.1	ng/ul	201677	3	14.251	1871760	20.0
Dibenzofuran, 4...	15.804	2.3	ng/ul	252229	4	17.069	2174860	20.0
Dibenzothiophene	16.898	3.8	ng/ul	411027	4	17.069	2174860	20.0
Phenanthrene, 2...	17.992	3.0	ng/ul	322570	4	17.069	2174860	20.0
1H-Indene, 2-ph...	18.051	4.1	ng/ul	449666	4	17.069	2174860	20.0
4H-Cyclopenta[d...	18.186	7.3	ng/ul	798670	4	17.069	2174860	20.0
Naphthalene, 2-...	18.510	2.2	ng/ul	237662	4	17.069	2174860	20.0
11H-Benzo[b]flu...	20.127	2.7	ng/ul	516096	5	21.516	3879880	20.0
11H-Benzo[a]flu...	20.221	3.0	ng/ul	576483	5	21.516	3879880	20.0
Benzo[e]pyrene	24.492	2.4	ng/ul	306845	6	24.792	2569190	20.0