LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072623\
Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4733DL

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-BP072523.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP016390.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.864	91	98	103	rBV3	92311	220879	2.19%	0.225%
2	3.981	113	118	122	rVV	107032	145443	1.44%	0.148%
3	4.175	145	151	161	rVB2	166852	298910	2.96%	0.305%
4	4.775	248	253	258	rVV	525187	790276	7.84%	0.806%
5	4.822	258	261	266	rVV	79310	114828	1.14%	0.117%
6	7.222	663	669	677	rBV2	153994	299506	2.97%	0.306%
7	7.428	695	704	710	rVB	124266	201152	1.99%	0.205%
8	7.611	729	735	741	rBV	135507	223546	2.22%	0.228%
9	8.093	810	817	827	rVB	1608811	2508762	24.88%	2.560%
10	8.787	928	935	941	rBV	163963	273746	2.71%	0.279%
11	8.934	951	960	968	rVB	67857	127048	1.26%	0.130%
12	9.281	1006	1019	1026	rBV	132529	229642	2.28%	0.234%
13	10.005	1133	1142	1147	rBV	74334	137977	1.37%	0.141%
14	10.528	1221	1231	1236	rBV	176562	326626	3.24%	0.333%
15	10.928	1287	1299	1305	rBV	2351221	3991954	39.59%	4.073%
16	14.051	1824	1830	1835	rVV2	76519	137650	1.37%	0.140%
17	14.146	1841	1846	1850	rVV	323683	471226	4.67%	0.481%
18	14.434	1887	1895	1898	rBV	376740	614006	6.09%	0.626%
19	14.463	1898	1900	1906	rVB	235519	292059	2.90%	0.298%
20	14.734	1935	1946	1953	rBV	3720077	5391975	53.47%	5.501%
21	14.798	1953	1957	1962	rVV	123333	196743	1.95%	0.201%
22	14.851	1962	1966	1972	rVV	87838	125535	1.24%	0.128%
23	14.916	1972	1977	1988	rVB2	89251	180550	1.79%	0.184%
24	15.128	2002	2013	2019	rBV2	85263	226460	2.25%	0.231%
25	15.728	2110	2115	2119	rVV	500908	706557	7.01%	0.721%
26	15.781	2119	2124	2131	rVB	368192	554085	5.50%	0.565%
27	15.987	2154	2159	2164	rVV	313013	447832	4.44%	0.457%
28	16.069	2169	2173	2179	rVV	116873	179072	1.78%	0.183%
29	16.216	2193	2198	2206	rVB2	110955	214874	2.13%	0.219%
30	16.428	2226	2234	2239	rBV8	41367	137893	1.37%	0.141%
31	16.710	2277	2282	2287	rBV	140742	195902	1.94%	0.200%
32	16.787	2287	2295	2299	rVV	191767	363810	3.61%	0.371%
33	16.834	2299	2303	2308	rVV	104451	143670	1.42%	0.147%
34	16.934	2316	2320	2325	rVB	102867	141357	1.40%	0.144%
35	17.010	2325	2333	2337	rBV2	145824	253907	2.52%	0.259%
36	17.051	2337	2340	2351	rVB3	69791	127509	1.26%	0.130%

LSC Area Percent Report

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Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4733DL

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-BP072523.MA.M
Title : SVOA CALIBRATION

37	17.210	2362	2367	2372	rBV3	89425	172543	1.71%	0.176%
38	17.316	2380	2385	2389	rBV	267052	361295	3.58%	0.369%
39	17.363	2389	2393	2396	rBV2	89170	128897	1.28%	0.132%
40	17.392	2396	2398	2405	rVB	114200	139463	1.38%	0.142%
41	17.498	2407	2416	2419	rBV	4470991	6531183	64.77%	6.664%
42	17.539	2419	2423	2429	rVB	4509507	6102907	60.53%	6.227%
43	17.598	2429	2433	2435	rBV2	457495	605405	6.00%	0.618%
44	17.634	2435	2439	2443	rVB	710508	911415	9.04%	0.930%
45	17.904	2481	2485	2488	rBV2	87268	126860	1.26%	0.129%
46	17.934	2488	2490	2498	rVB5	89948	196803	1.95%	0.201%
47	18.010	2498	2503	2508	rBV	150580	217454	2.16%	0.222%
48	18.069	2508	2513	2521	rVB3	344971	596669	5.92%	0.609%
49	18.204	2530	2536	2543	rVB	135878	281991	2.80%	0.288%
50	18.351	2552	2561	2565	rBV	896920	1340495	13.29%	1.368%
51	18.398	2565	2569	2577	rVB	1032588	1397094	13.86%	1.425%
52	18.475	2578	2582	2587	rBV	314670	403583	4.00%	0.412%
53	18.539	2587	2593	2597	rBV3	1200211	2293076	22.74%	2.340%
54	18.575	2597	2599	2604	rVV	696065	806098	7.99%	0.822%
55	18.792	2630	2636	2639	rBV6	77769	151971	1.51%	0.155%
56	18.834	2639	2643	2648	rVV	539930	751385	7.45%	0.767%
57	18.881	2648	2651	2657	rVB2	150468	203381	2.02%	0.208%
58	19.063	2674	2682	2686	rBV4	148838	312594	3.10%	0.319%
59	19.110	2687	2690	2693	rVV	158271	194248	1.93%	0.198%
60	19.151	2693	2697	2701	rVB2	125869	163534	1.62%	0.167%
61	19.228	2705	2710	2714	rVV	411695	631307	6.26%	0.644%
62	19.292	2714	2721	2724	rVV2	256159	608818	6.04%	0.621%
63	19.322	2724	2726	2729	rVV	225551	241906	2.40%	0.247%
64	19.375	2729	2735	2747	rVV8	167006	582652	5.78%	0.594%
65	19.492	2750	2755	2760	rVV	5114638	6582853	65.29%	6.716%
66	19.545	2762	2764	2768	rVV	129684	215063	2.13%	0.219%
67	19.581	2768	2770	2775	rVV2	146271	223483	2.22%	0.228%
68	19.639	2776	2780	2784	rVV	337873	451845	4.48%	0.461%
69	19.733	2791	2796	2803	rVV2	161836	339828	3.37%	0.347%
70	19.816	2804	2810	2812	rVV2	1077156	1664516	16.51%	1.698%
71	19.851	2812	2816	2822	rVV	4710890	6390169	63.37%	6.520%
72	19.910	2822	2826	2830	rVB	247525	326309	3.24%	0.333%
73	20.016	2841	2844	2848	rVB	210653	241947	2.40%	0.247%
74	20.157	2862	2868	2875	rBV3	359837	879353	8.72%	0.897%
75	20.233	2877	2881	2885	rVV3	181239	246359	2.44%	0.251%
76	20.322	2893	2896	2904	rVB	922169	1267571	12.57%	1.293%

LSC Area Percent Report

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Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

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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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Title : SVOA CALIBRATION

77	20.422	2909	2913	2917	rBV	886046	1073973	10.65%	1.096%
78	20.480	2917	2923	2926	rVB2	457254	684191	6.79%	0.698%
79	20.616	2942	2946	2950	rBV	359872	435149	4.32%	0.444%
80	20.663	2950	2954	2959	rVB	421254	545540	5.41%	0.557%
81	20.751	2966	2969	2978	rVB3	176316	292028	2.90%	0.298%
82	21.010	3010	3013	3017	rBV5	102711	173966	1.73%	0.177%
83	21.222	3046	3049	3052	rBV	276442	306276	3.04%	0.312%
84	21.257	3052	3055	3059	rVV	414170	490928	4.87%	0.501%
85	21.298	3059	3062	3067	rVB2	251843	340469	3.38%	0.347%
86	21.463	3087	3090	3099	rVB2	203440	313101	3.11%	0.319%
87	21.586	3103	3111	3115	rBV2	5236627	10083194	100.00%	10.288%
88	21.627	3115	3118	3123	rVB	1797157	2331774	23.13%	2.379%
89	21.751	3136	3139	3145	rBV2	135460	199697	1.98%	0.204%
90	21.810	3145	3149	3153	rBV	241588	339096	3.36%	0.346%
91	22.192	3210	3214	3219	rVB2	355046	527353	5.23%	0.538%
92	22.251	3221	3224	3229	rVB	201782	271662	2.69%	0.277%
93	22.416	3248	3252	3255	rBV3	193036	291471	2.89%	0.297%
94	22.457	3256	3259	3265	rVB2	238835	318139	3.16%	0.325%
95	23.374	3407	3415	3421	rBV2	961875	2873916	28.50%	2.932%
96	23.574	3446	3449	3458	rVB	206451	360615	3.58%	0.368%
97	23.945	3507	3512	3518	rBV	516235	993899	9.86%	1.014%
98	24.004	3519	3522	3526	rVV2	242326	399649	3.96%	0.408%
99	24.063	3526	3532	3542	rVB	1037425	2234731	22.16%	2.280%
100	24.180	3545	3552	3559	rBV	2668017	5852697	58.04%	5.971%

Sum of corrected areas: 98010804

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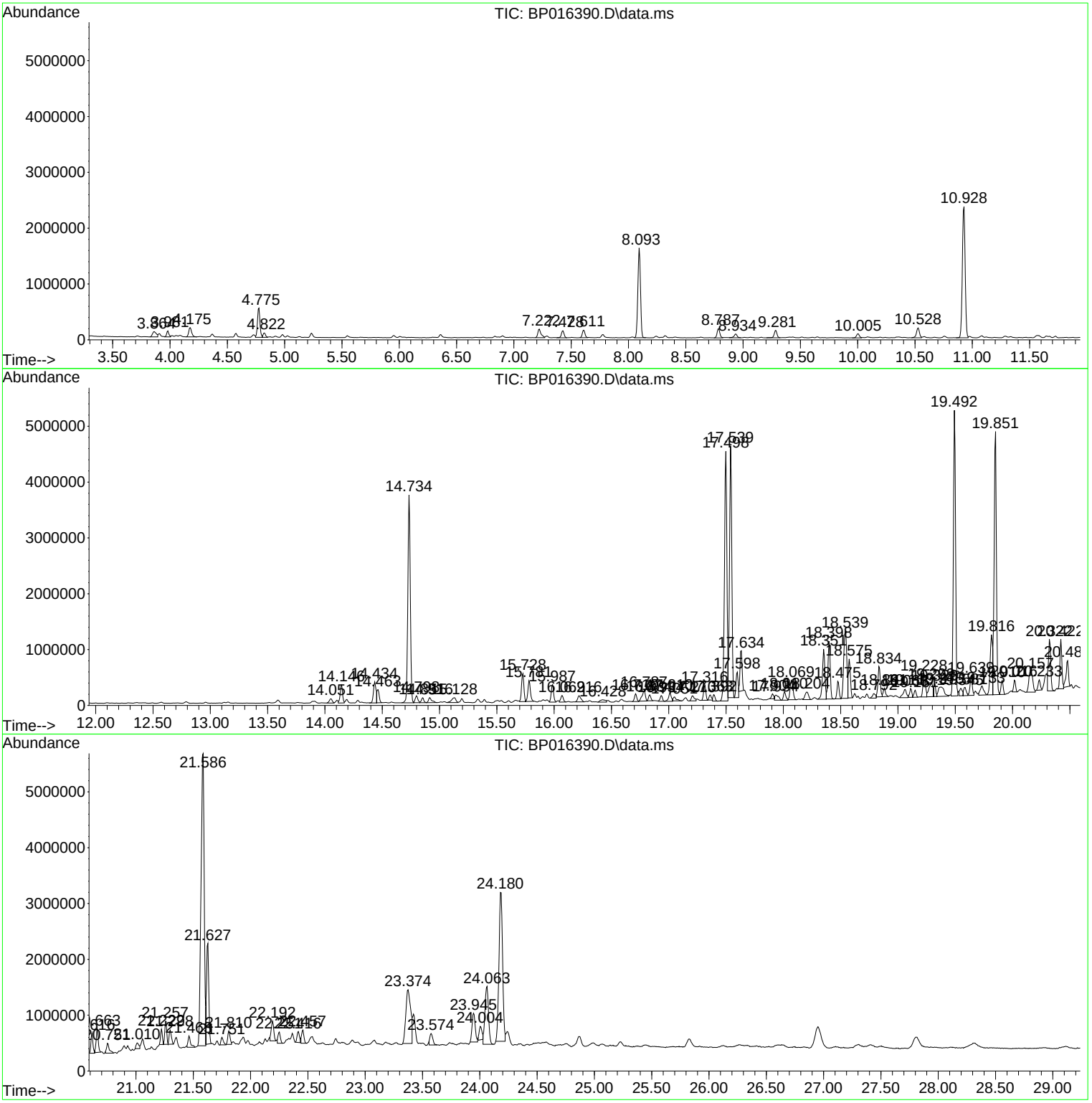
Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.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\BP072623\
Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
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Sample : 03534-17DL 5X
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Instrument :
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ClientSampleId :
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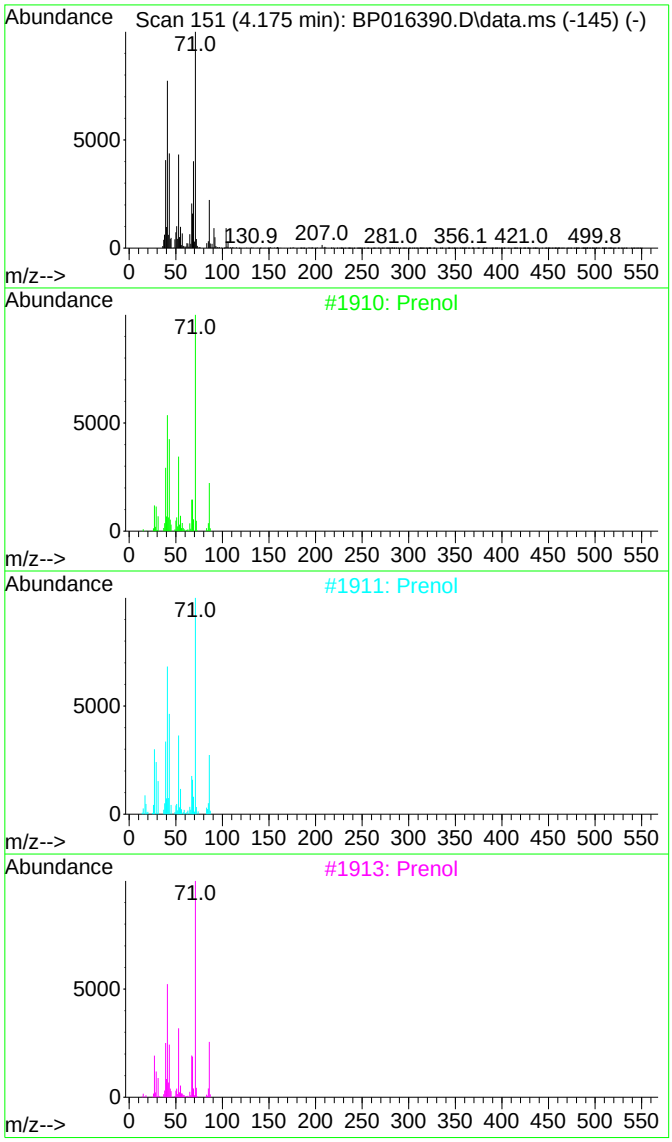
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

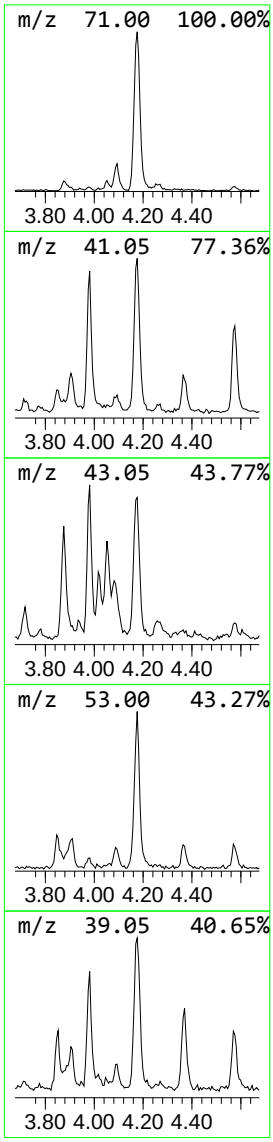
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Prenol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.175	2.38 ng/ul	298910	1,4-Dichlorobenzene-d4	8.093

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Prenol	86	C5H10O	000556-82-1	76
2	Prenol	86	C5H10O	000556-82-1	72
3	Prenol	86	C5H10O	000556-82-1	38
4	2-Butene, 2-chloro-3-methyl-	104	C5H9Cl	017773-65-8	37
5	1-Butene, 2-chloro-3-methyl-	104	C5H9Cl	017773-64-7	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072623\
Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

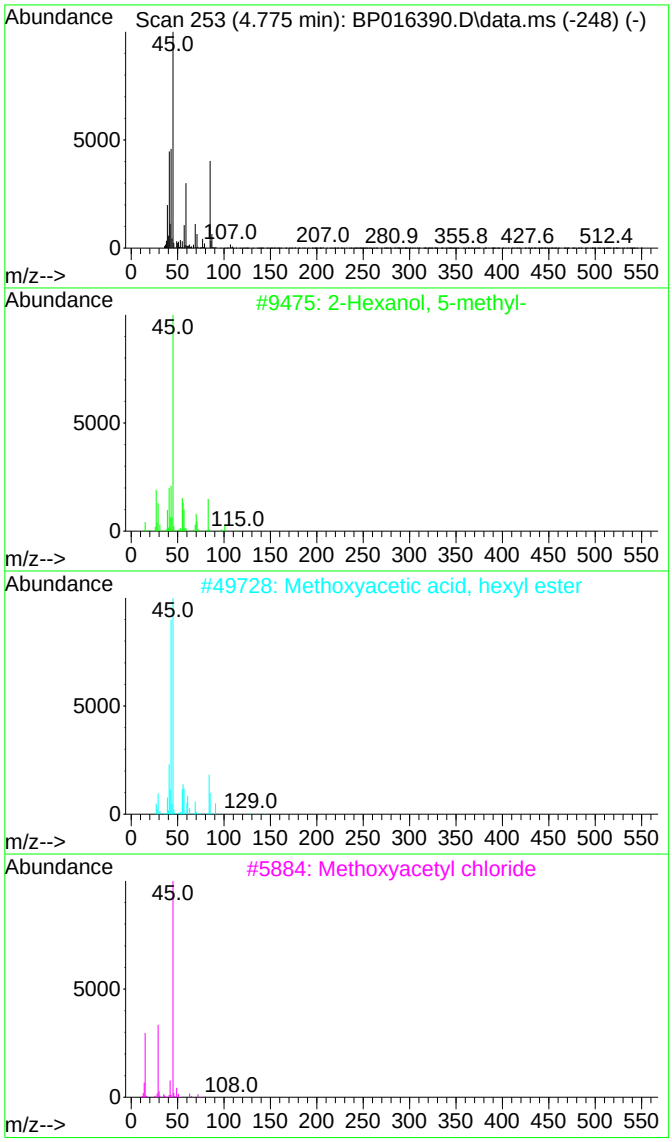
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

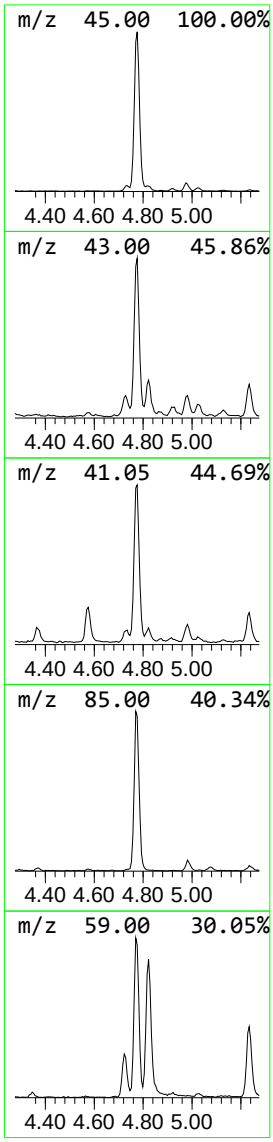
****

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.775	6.30 ng/ul	790276	1,4-Dichlorobenzene-d4	8.093

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 5-methyl-	116	C7H16O	000627-59-8	37
2	Methoxyacetic acid, hexyl ester	174	C9H18O3	145747-16-6	33
3	Methoxyacetyl chloride	108	C3H5ClO2	038870-89-2	32
4	5-Hexen-2-ol	100	C6H12O	000626-94-8	32
5	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072623\
Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
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Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

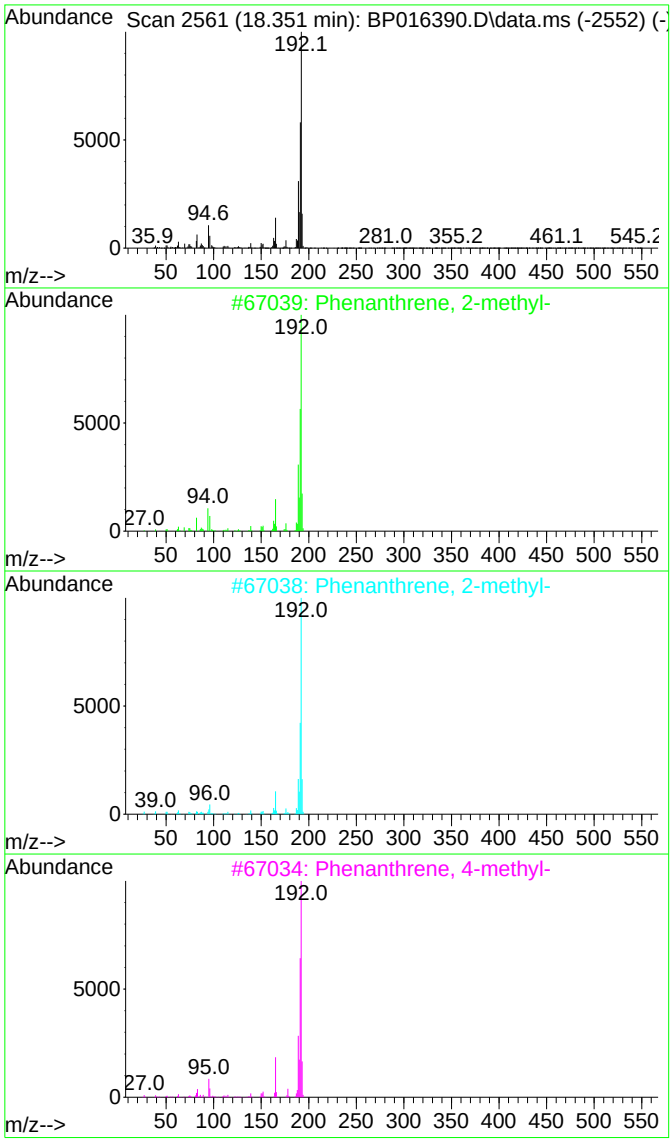
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

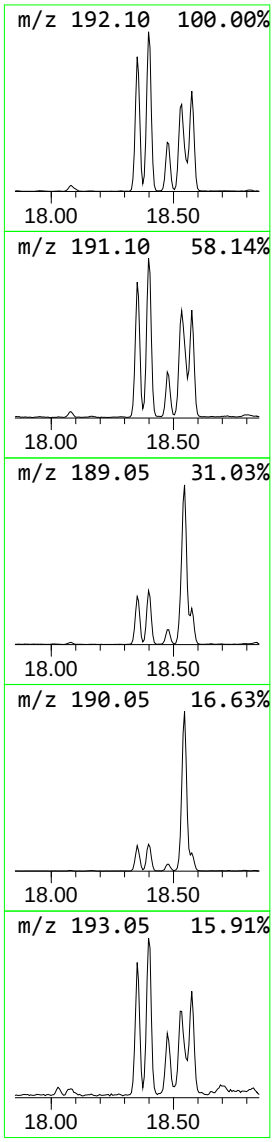
****

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.351	4.10 ng/ul	1340500	Phenanthrene-d10	17.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94





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Acq On : 26 Jul 2023 23:47
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Misc :
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Instrument :
BNA_P
ClientSampleId :
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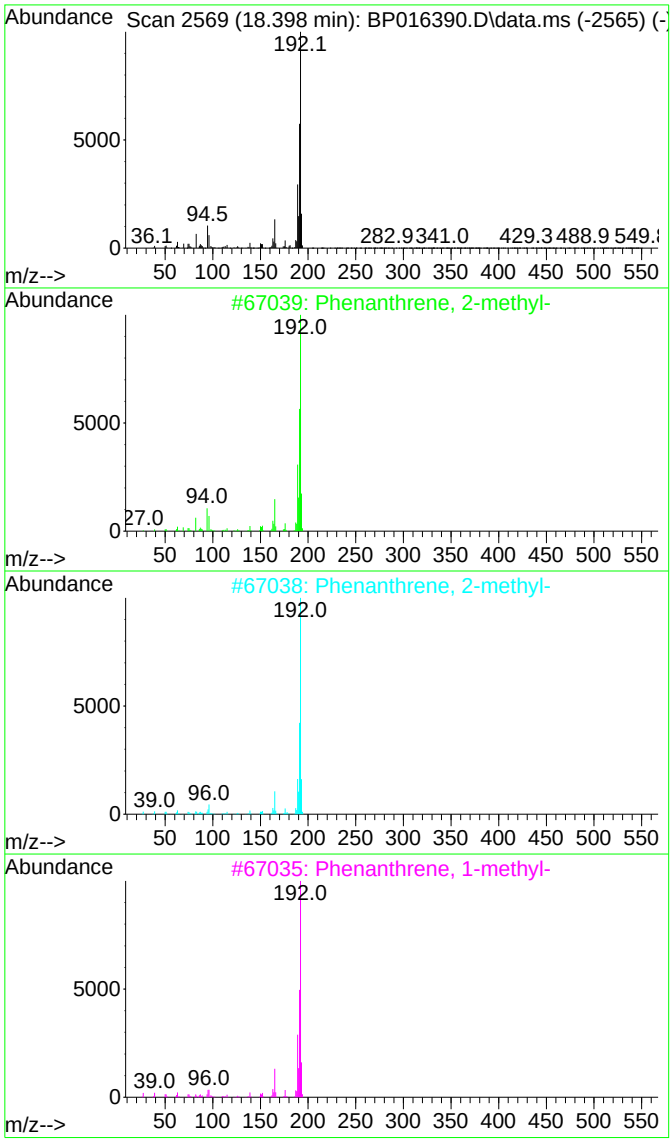
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

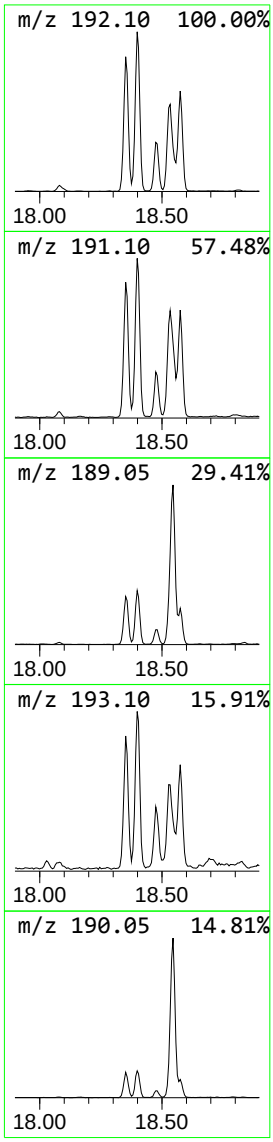
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	4.28 ng/ul	1397090	Phenanthrene-d10	17.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072623\
Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

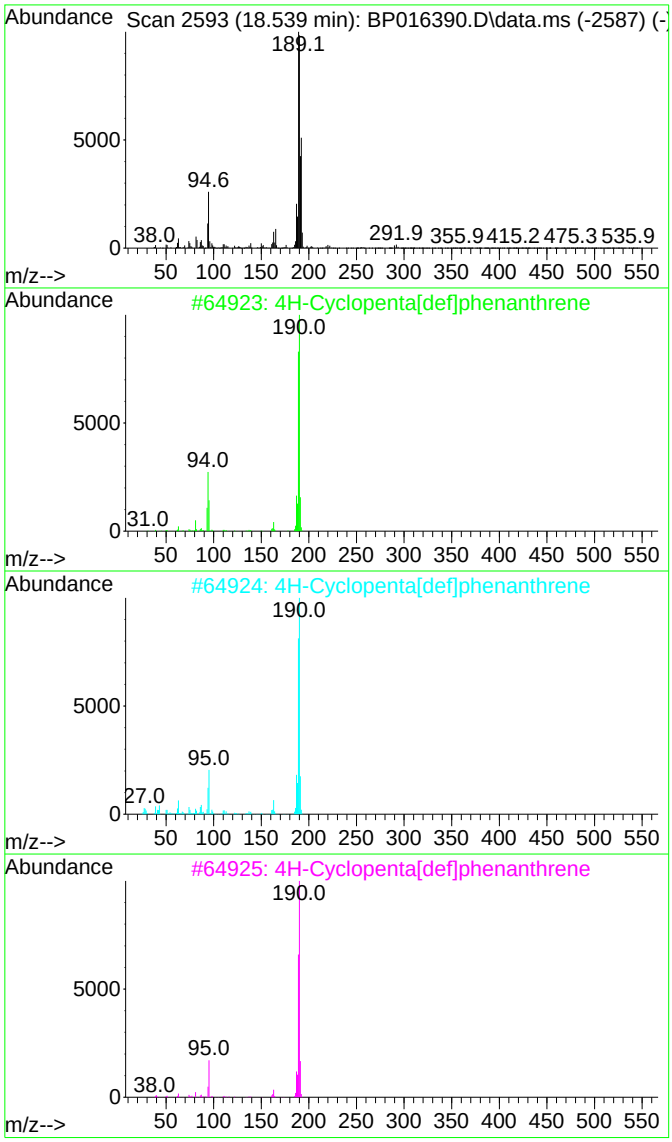
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

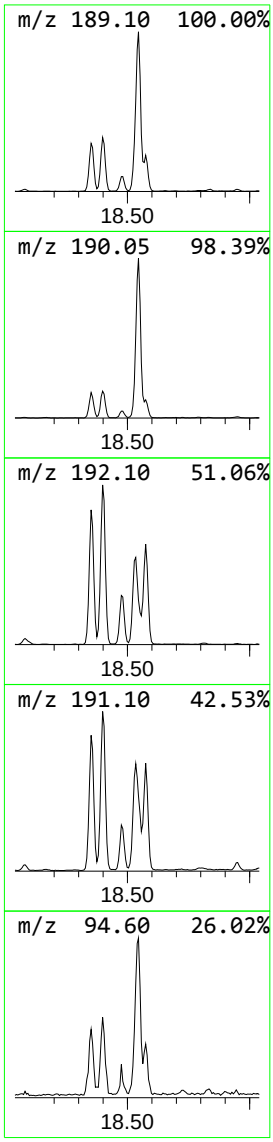
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	7.02 ng/ul	2293080	Phenanthrene-d10	17.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
4	4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	38
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18





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Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

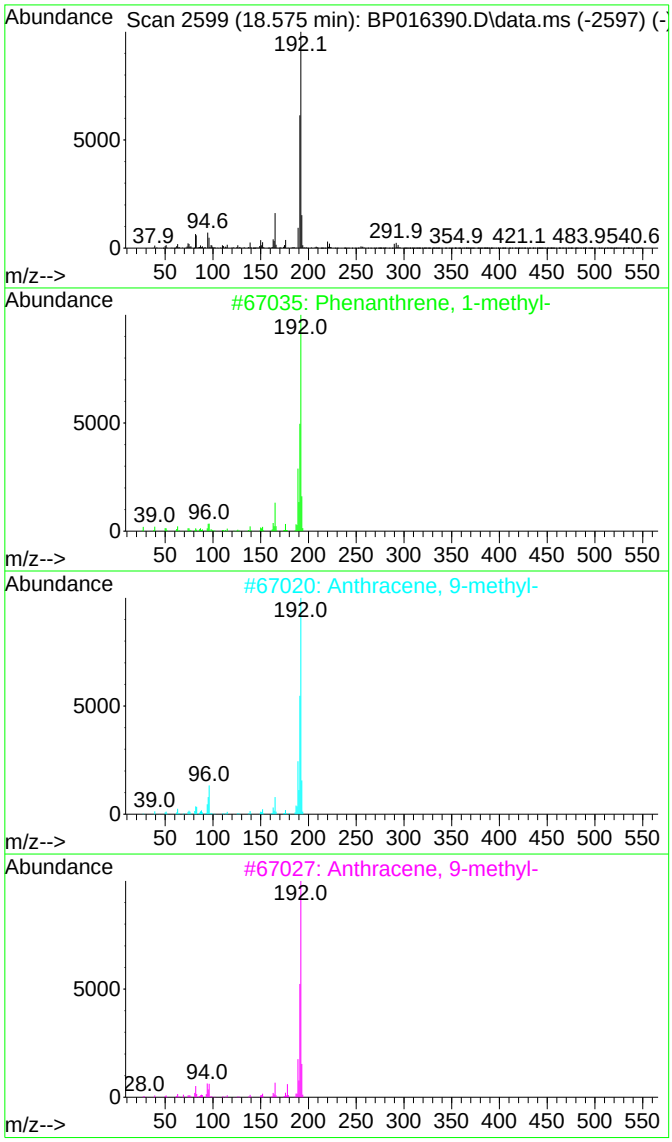
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

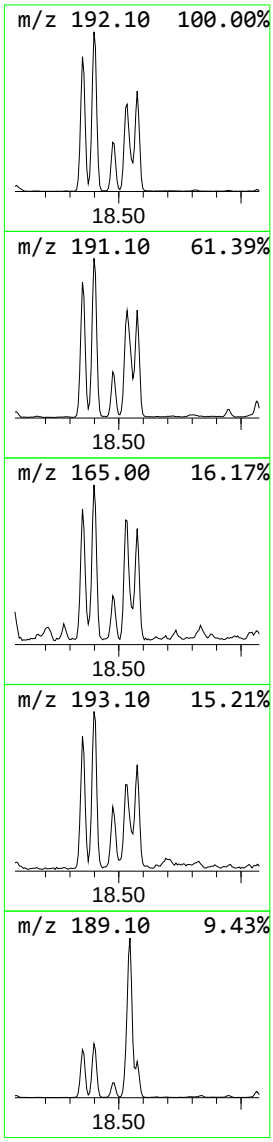
****

Peak Number 6 Anthracene, 9-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.575	2.47 ng/ul	806098	Phenanthrene-d10	17.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87
5	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	87





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Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

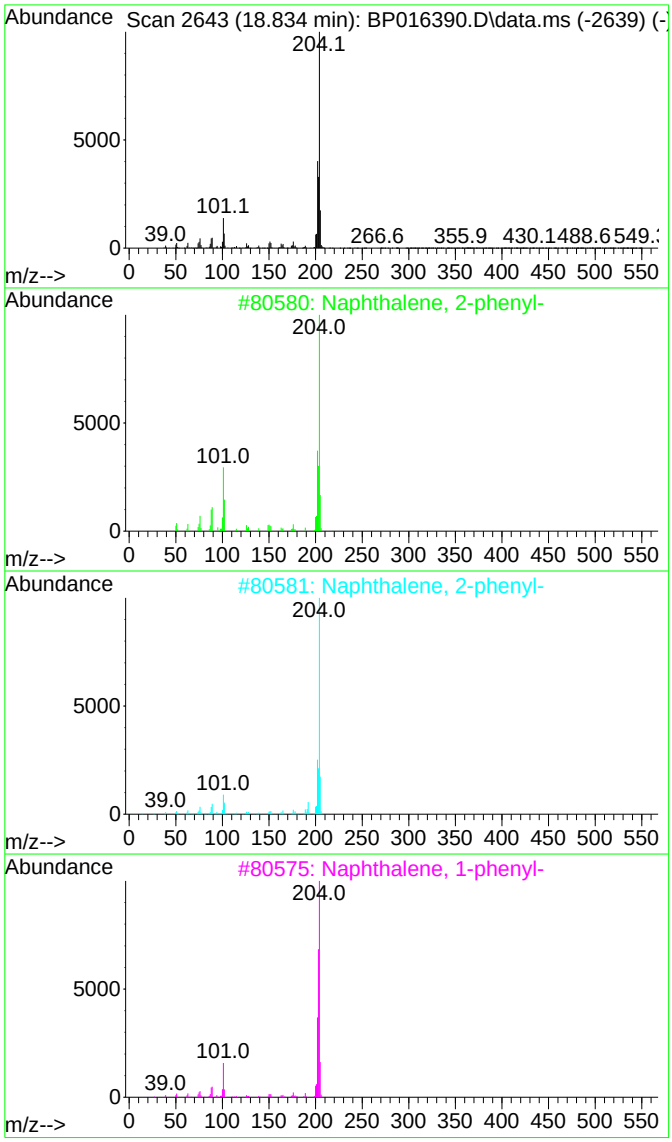
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

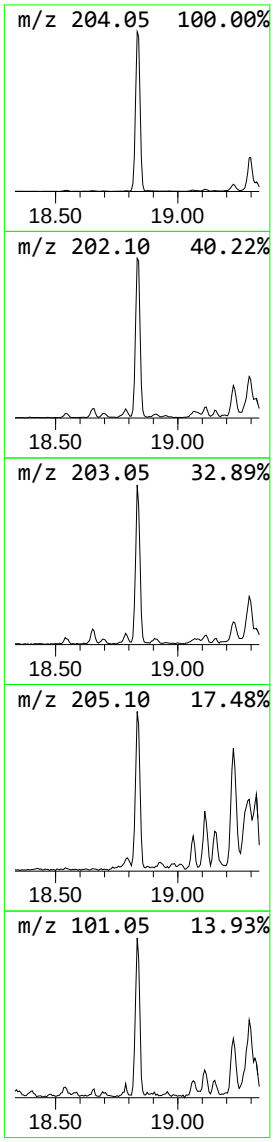
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.834	2.30 ng/ul	751385	Phenanthrene-d10	17.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	95
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
3	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	83
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	68





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Data File : BP016390.D
Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

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

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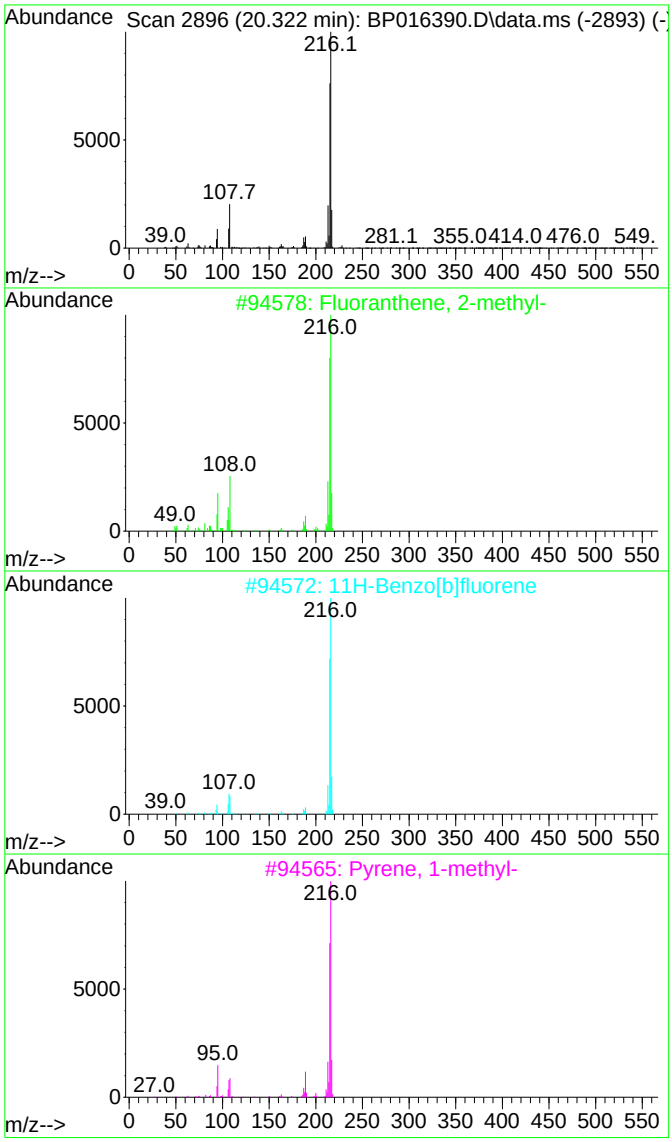
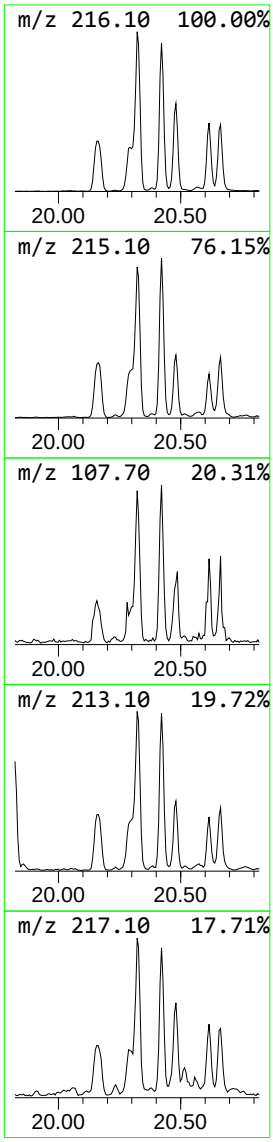
Peak Number 8 Fluoranthene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.322	2.51 ng/ul	1267570	Chrysene-d12	21.586

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	90





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Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

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

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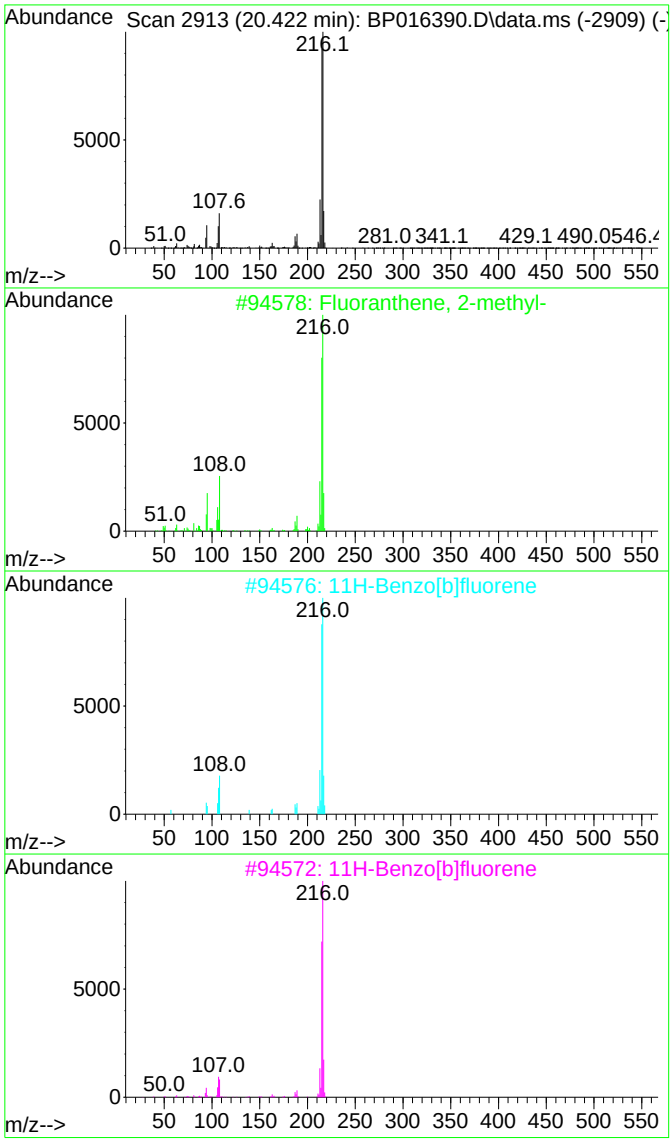
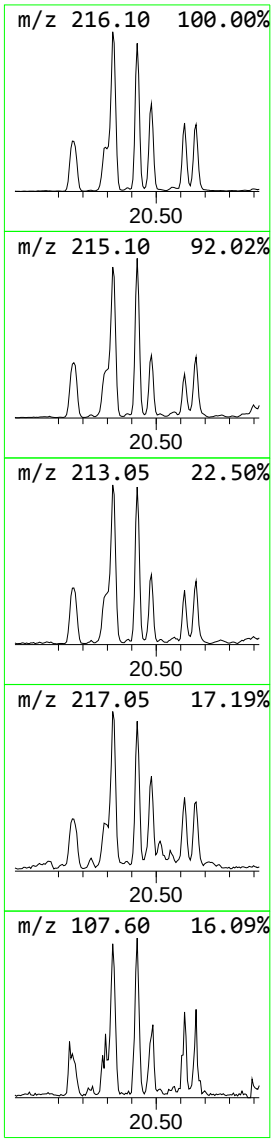
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.422	2.13 ng/ul	1073970	Chrysene-d12	21.586

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87





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Acq On : 26 Jul 2023 23:47
Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4733DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.MA.M
Quant Title : SVOA CALIBRATION

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

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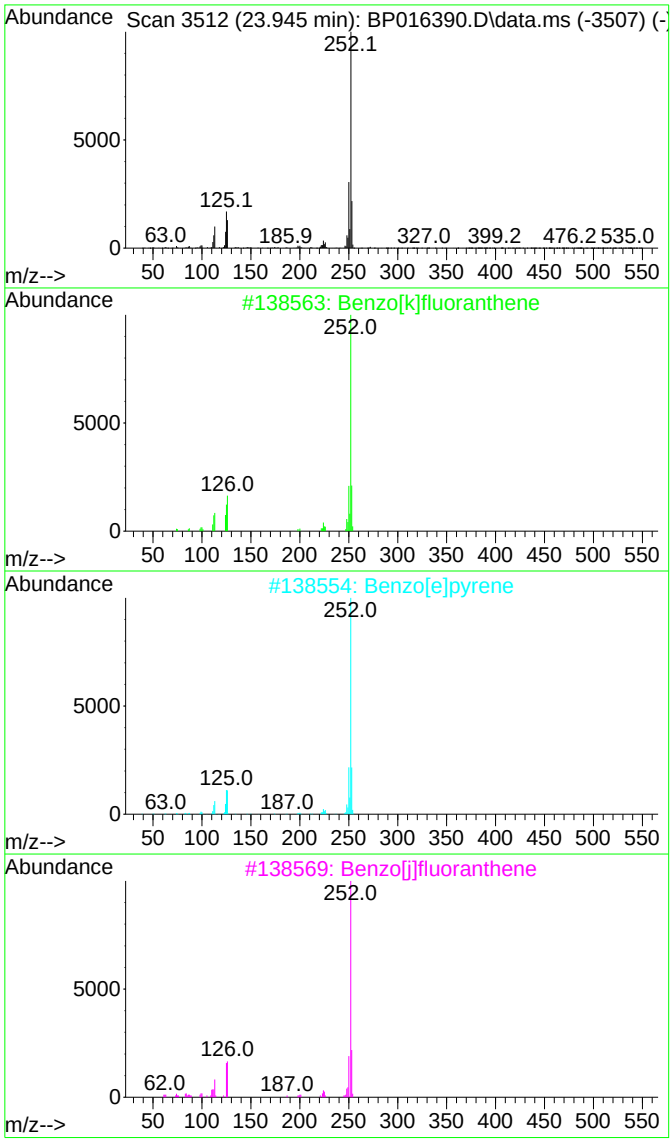
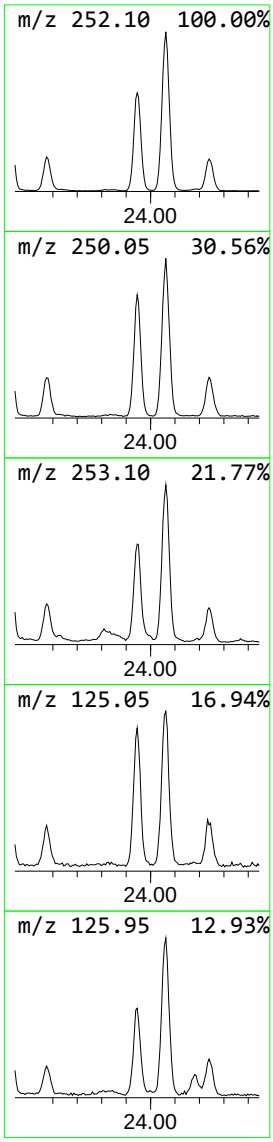
Peak Number 10 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.945	3.40 ng/ul	993899	Perylene-d12	24.186

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
2	Benzo[e]pyrene	252	C20H12	000192-97-2	96
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	96
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5	Benzo[a]pyrene	252	C20H12	000050-32-8	93





Tentatively Identified Compound (LSC) summary

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Data File : BP016390.D
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Operator : MA/JU
Sample : 03534-17DL 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072523.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
Prenol	4.175	2.4	ng/ul	298910	1	8.093	2508760	20.0
unknown-01	4.775	6.3	ng/ul	790276	1	8.093	2508760	20.0
Phenanthrene, 2...	18.351	4.1	ng/ul	1340500	4	17.498	6531180	20.0
Phenanthrene, 1...	18.398	4.3	ng/ul	1397090	4	17.498	6531180	20.0
4H-Cyclopenta[d...	18.539	7.0	ng/ul	2293080	4	17.498	6531180	20.0
Anthracene, 9-m...	18.575	2.5	ng/ul	806098	4	17.498	6531180	20.0
Naphthalene, 2-...	18.834	2.3	ng/ul	751385	4	17.498	6531180	20.0
Fluoranthene, 2...	20.322	2.5	ng/ul	1267570	5	21.586	10083200	20.0
11H-Benzo[b]flu...	20.422	2.1	ng/ul	1073970	5	21.586	10083200	20.0
Benzo[e]pyrene	23.945	3.4	ng/ul	993899	6	24.186	5852700	20.0