LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

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-BP042122.M
Title : SVOA CALIBRATION

Signal : TIC: BP010062.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.799	118	121	133	rBV2	32250	65344	1.26%	0.134%
2	3.893	133	137	144	rVB	28707	40370	0.78%	0.083%
3	4.117	172	175	183	rVB3	9753	17225	0.33%	0.035%
4	7.099	674	682	692	rBV	104191	179280	3.47%	0.367%
5	7.264	704	710	721	rBV	83265	142367	2.75%	0.291%
6	7.469	735	745	754	rBV	110738	184993	3.58%	0.378%
7	7.940	817	825	841	rBV	648173	1082257	20.93%	2.213%
8	8.646	938	945	953	rBV	131034	233732	4.52%	0.478%
9	9.099	1014	1022	1032	rBV	110480	196541	3.80%	0.402%
10	9.546	1093	1098	1103	rVB6	7439	13103	0.25%	0.027%
11	9.822	1138	1145	1152	rBV2	87200	154746	2.99%	0.316%
12	10.363	1229	1237	1245	rBV	147006	260218	5.03%	0.532%
13	10.446	1249	1251	1260	rVB10	4773	11616	0.22%	0.024%
14	10.746	1291	1302	1317	rBV	1016352	1753536	33.92%	3.586%
15	10.881	1317	1325	1331	rBV2	35149	72841	1.41%	0.149%
16	13.893	1831	1837	1842	rBV5	11688	21547	0.42%	0.044%
17	13.975	1842	1851	1857	rBV	307505	429084	8.30%	0.877%
18	14.263	1889	1900	1912	rBV	309659	535305	10.35%	1.095%
19	14.569	1941	1952	1959	rBV	1678212	2511317	48.58%	5.136%
20	14.634	1959	1963	1971	rVB2	25231	46768	0.90%	0.096%
21	14.763	1979	1985	2000	rBV	66130	137411	2.66%	0.281%
22	14.975	2015	2021	2027	rVV2	18892	33476	0.65%	0.068%
23	15.045	2028	2033	2037	rVB5	9513	17863	0.35%	0.037%
24	15.193	2052	2058	2063	rBV10	9302	19604	0.38%	0.040%
25	15.363	2080	2087	2096	rBV10	8053	26734	0.52%	0.055%
26	15.563	2114	2121	2126	rBV	435302	656763	12.70%	1.343%
27	15.616	2126	2130	2136	rVV	41440	67632	1.31%	0.138%
28	15.675	2136	2140	2148	rVV2	65378	99228	1.92%	0.203%
29	15.745	2148	2152	2155	rVB6	10602	15546	0.30%	0.032%
30	15.934	2177	2184	2190	rVB8	15108	39711	0.77%	0.081%
31	16.010	2193	2197	2200	rBV	26564	37426	0.72%	0.077%
32	16.045	2200	2203	2206	rVV4	11146	19682	0.38%	0.040%
33	16.087	2206	2210	2215	rVB2	16288	28197	0.55%	0.058%
34	16.210	2227	2231	2233	rBV5	7955	11506	0.22%	0.024%
35	16.310	2239	2248	2253	rBV	110500	174513	3.38%	0.357%
36	16.351	2253	2255	2259	rVB4	10119	12718	0.25%	0.026%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFFE0DL

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-BP042122.M
Title : SVOA CALIBRATION

37	16.404	2259	2264	2268	rBV	298812	387242	7.49%	0.792%
38	16.463	2268	2274	2280	rVV2	280178	568396	10.99%	1.162%
39	16.522	2280	2284	2288	rVV2	326312	444626	8.60%	0.909%
40	16.569	2288	2292	2296	rVV	173382	246392	4.77%	0.504%
41	16.622	2296	2301	2302	rVV2	76565	113076	2.19%	0.231%
42	16.645	2302	2305	2307	rVV	97016	143275	2.77%	0.293%
43	16.669	2307	2309	2313	rVV	118516	148365	2.87%	0.303%
44	16.722	2313	2318	2325	rVV3	228047	461491	8.93%	0.944%
45	16.792	2325	2330	2333	rVV	299613	421581	8.15%	0.862%
46	16.828	2333	2336	2339	rVV2	88135	140345	2.71%	0.287%
47	16.863	2339	2342	2351	rVB2	163882	248567	4.81%	0.508%
48	16.975	2356	2361	2365	rBV	44030	69289	1.34%	0.142%
49	17.139	2381	2389	2397	rVB3	48443	109319	2.11%	0.224%
50	17.216	2397	2402	2403	rBV3	12141	17593	0.34%	0.036%
51	17.322	2413	2420	2424	rBV	2019503	3017133	58.36%	6.170%
52	17.363	2424	2427	2432	rVB	975490	1282941	24.82%	2.624%
53	17.422	2432	2437	2441	rBV	465725	671299	12.99%	1.373%
54	17.516	2448	2453	2460	rBV2	65327	102883	1.99%	0.210%
55	17.722	2481	2488	2493	rBV2	115261	192311	3.72%	0.393%
56	17.839	2504	2508	2514	rBV4	19038	31174	0.60%	0.064%
57	17.898	2515	2518	2527	rVB2	31355	45866	0.89%	0.094%
58	17.986	2531	2533	2537	rVV5	15163	20945	0.41%	0.043%
59	18.039	2539	2542	2547	rVB2	12853	19234	0.37%	0.039%
60	18.186	2563	2567	2571	rBV	132111	201853	3.90%	0.413%
61	18.234	2571	2575	2580	rVB	186601	253348	4.90%	0.518%
62	18.316	2584	2589	2593	rBV4	20214	34214	0.66%	0.070%
63	18.375	2593	2599	2603	rBV2	196765	361708	7.00%	0.740%
64	18.410	2603	2605	2612	rVB	106720	121068	2.34%	0.248%
65	18.522	2621	2624	2628	rBV4	16749	20790	0.40%	0.043%
66	18.669	2644	2649	2651	rBV	120426	151196	2.92%	0.309%
67	18.704	2651	2655	2661	rVB	277659	394154	7.62%	0.806%
68	18.910	2686	2690	2694	rBV	36966	45191	0.87%	0.092%
69	18.957	2695	2698	2701	rVV	34510	40135	0.78%	0.082%
70	18.992	2701	2704	2708	rVB	32502	36752	0.71%	0.075%
71	19.075	2713	2718	2722	rBV2	119050	174057	3.37%	0.356%
72	19.204	2736	2740	2750	rBV	125540	216222	4.18%	0.442%
73	19.328	2756	2761	2767	rVB	2723784	3564098	68.94%	7.288%
74	19.422	2774	2777	2782	rVB2	49802	64770	1.25%	0.132%
75	19.516	2789	2793	2796	rBV2	50722	78203	1.51%	0.160%
76	19.563	2798	2801	2810	rVB	82207	154892	3.00%	0.317%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

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-BP042122.M
Title : SVOA CALIBRATION

77	19.651	2811	2816	2818	rBV3	1026569	1474273	28.52%	3.015%
78	19.681	2818	2821	2828	rVV2	2627634	3438813	66.52%	7.032%
79	19.745	2829	2832	2838	rVB2	98184	154405	2.99%	0.316%
80	19.851	2847	2850	2856	rBV	95813	123001	2.38%	0.252%
81	19.910	2857	2860	2865	rVB	104278	150643	2.91%	0.308%
82	19.998	2869	2875	2880	rBV3	129120	318690	6.16%	0.652%
83	20.163	2899	2903	2908	rVB	241289	343110	6.64%	0.702%
84	20.257	2915	2919	2923	rBV	173821	229208	4.43%	0.469%
85	20.316	2925	2929	2932	rVB2	158195	216653	4.19%	0.443%
86	20.451	2949	2952	2956	rBV	103151	132343	2.56%	0.271%
87	20.498	2957	2960	2963	rVB2	102147	127282	2.46%	0.260%
88	20.892	3023	3027	3033	rVB	255171	330439	6.39%	0.676%
89	21.057	3052	3055	3058	rVB	192275	229729	4.44%	0.470%
90	21.092	3058	3061	3063	rBV	132314	170933	3.31%	0.350%
91	21.180	3073	3076	3083	rVB2	257260	337809	6.53%	0.691%
92	21.404	3107	3114	3118	rBV2	2738597	5169662	100.00%	10.572%
93	21.445	3118	3121	3132	rVB	1313147	1963706	37.99%	4.016%
94	21.569	3139	3142	3146	rVB2	148590	191710	3.71%	0.392%
95	21.733	3167	3170	3188	rVB6	229493	640815	12.40%	1.310%
96	23.110	3398	3404	3410	rBV	1105206	2632346	50.92%	5.383%
97	23.639	3490	3494	3500	rBV	441886	815450	15.77%	1.668%
98	23.751	3508	3513	3523	rVB2	762907	1627142	31.47%	3.327%
99	23.857	3525	3531	3538	rBV2	1171241	2486179	48.09%	5.084%
100	26.415	3959	3966	3980	rVB3	439017	1430403	27.67%	2.925%

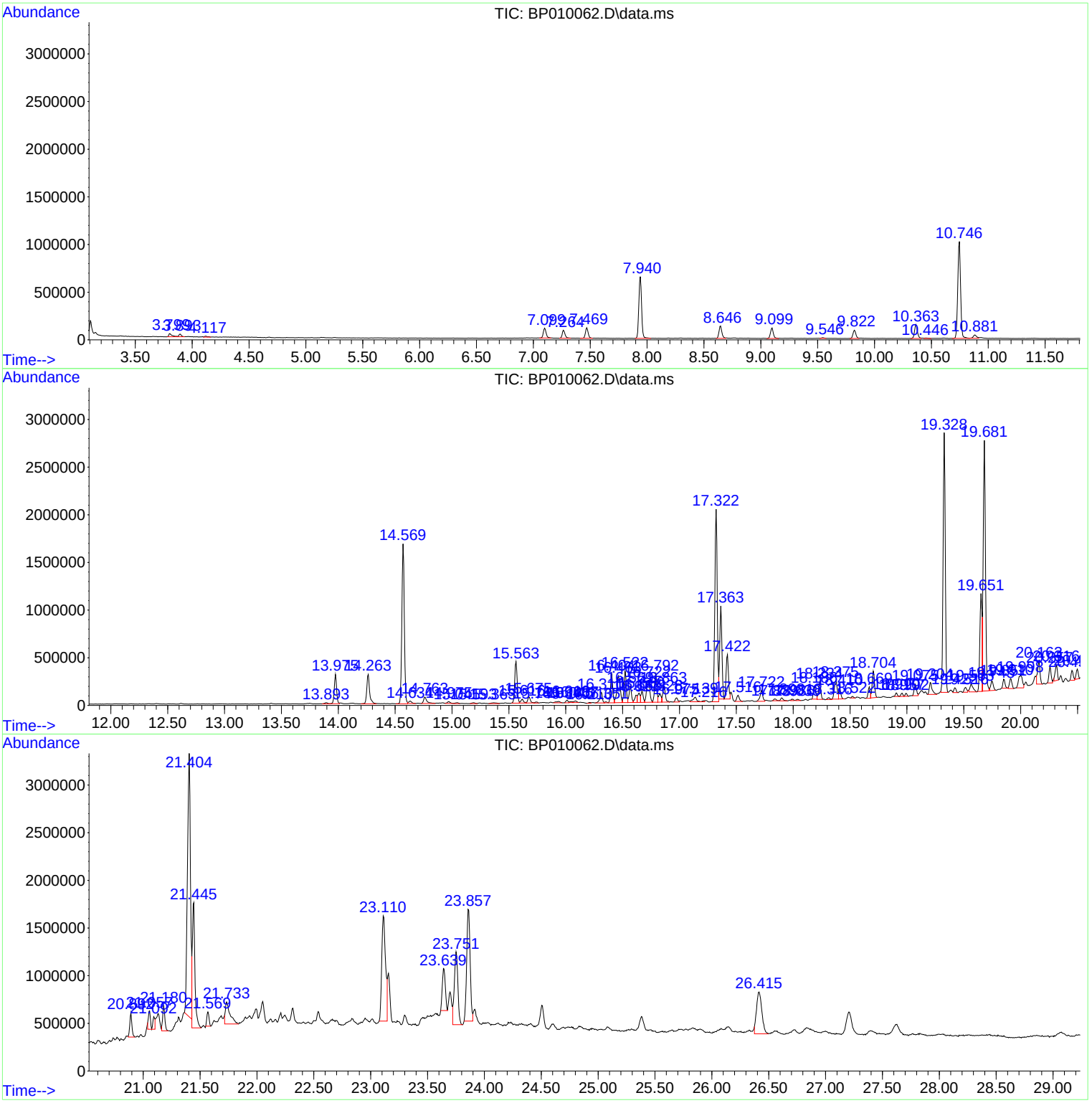
Sum of corrected areas: 48900938

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

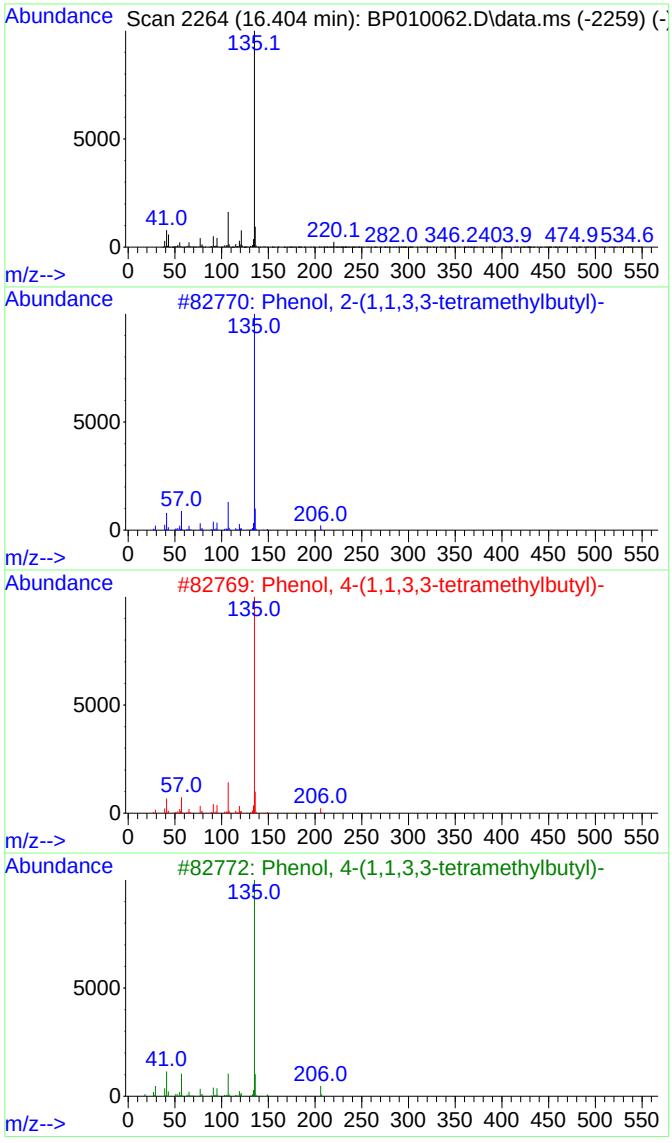
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

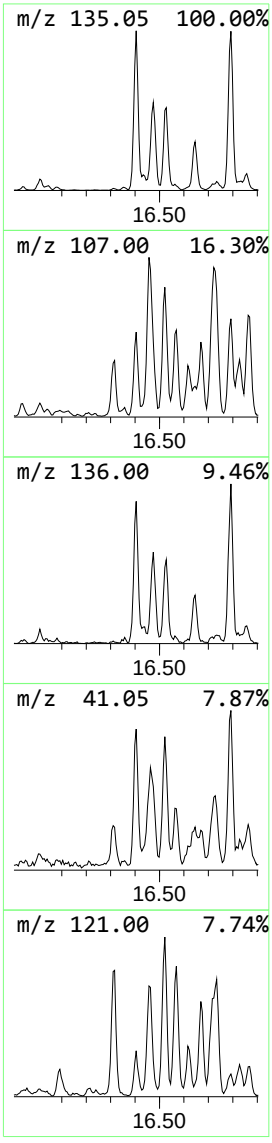
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Phenol, 2-(1,1,3,3-tetramet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.404	2.57 ng/ul	387242	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	86
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
3	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
4	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

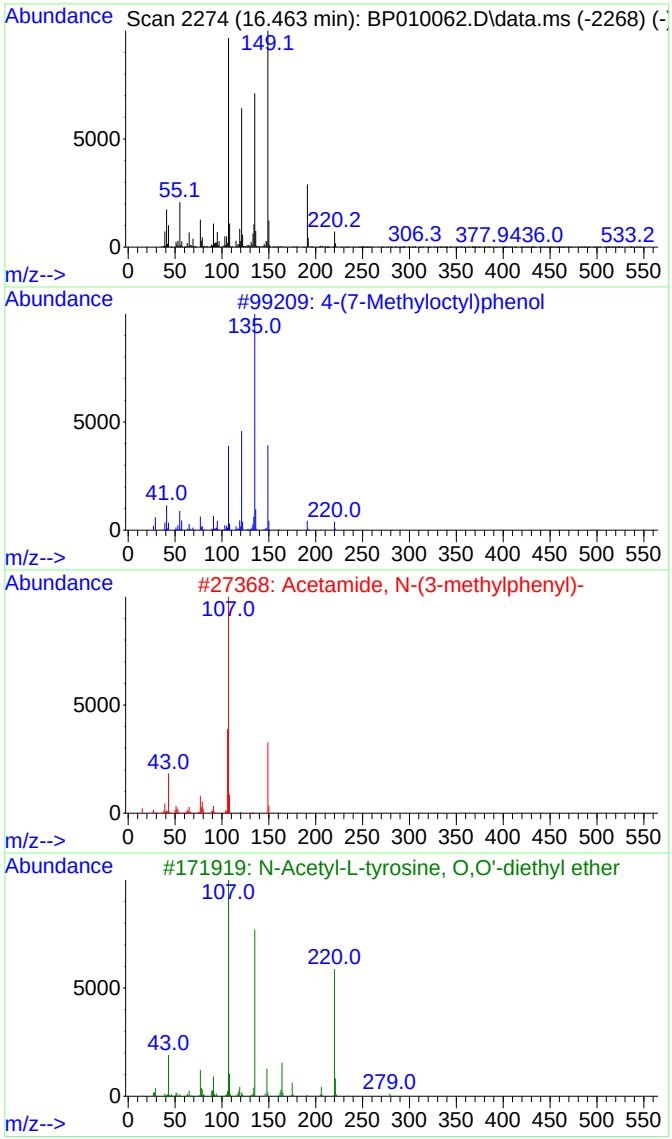
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

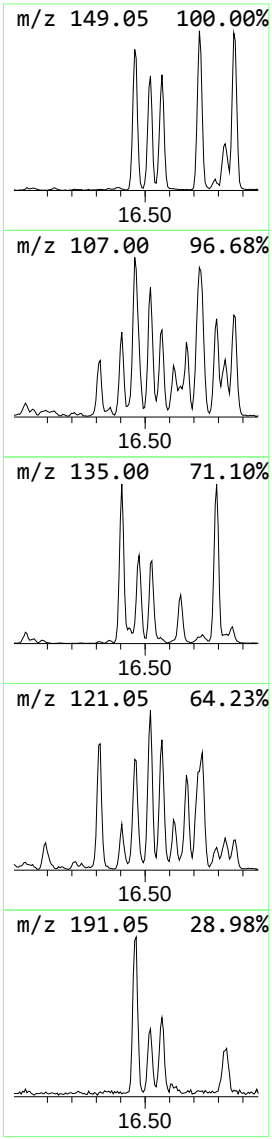
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 4-(7-Methyloctyl)phenol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.463	3.77 ng/ul	568396	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	62
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	52
3	N-Acetyl-L-tyrosine, O,O'-diethy...	279	C15H21NO4	1000453-79-9	38
4	Ethanol, 2-(3,3-dimethylbicyclo[...	166	C11H18O	002226-05-3	38
5	Phenol, 4-(2-methylpropyl)-	150	C10H14O	004167-74-2	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

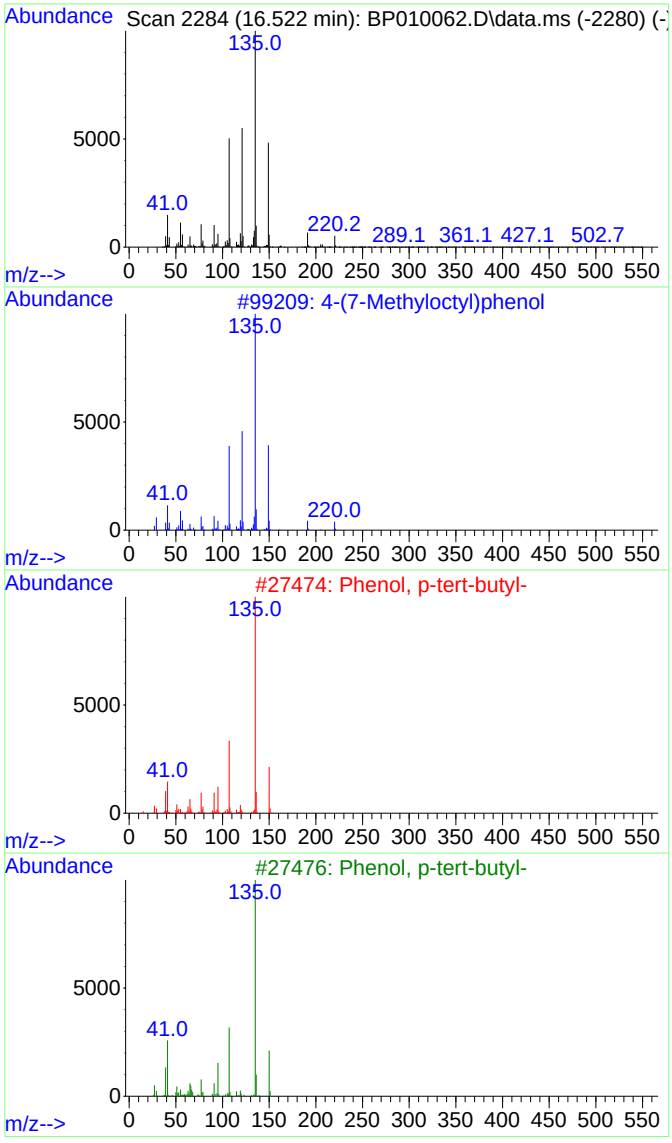
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

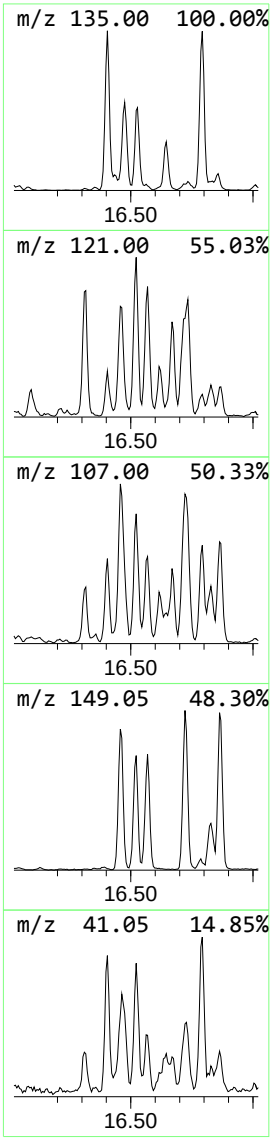
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.522	2.95 ng/ul	444626	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	98
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	58
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	52
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

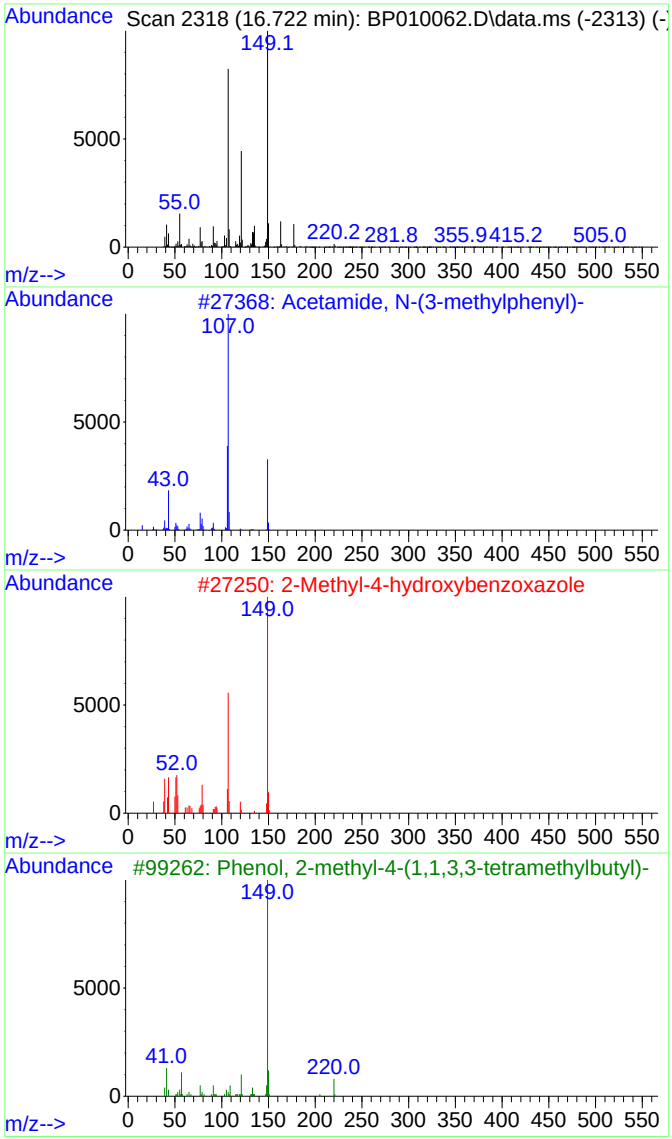
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

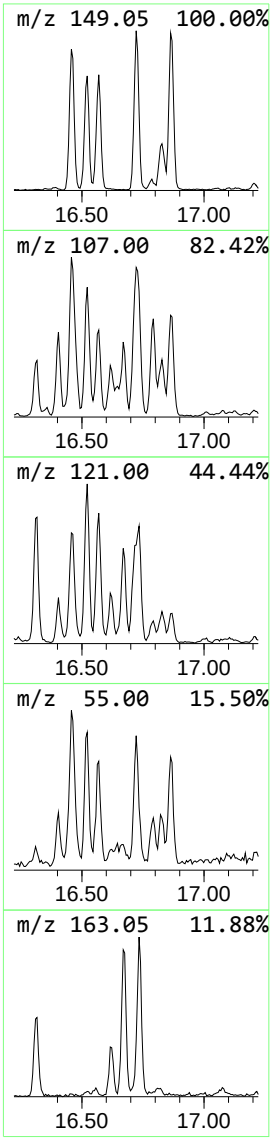
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Acetamide, N-(3-methylphenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.722	3.06 ng/ul	461491	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	59
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	45
4	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
5	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

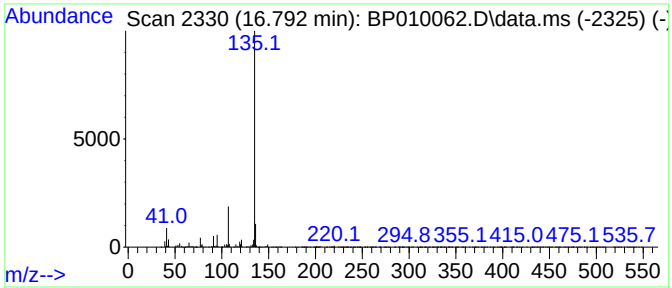
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

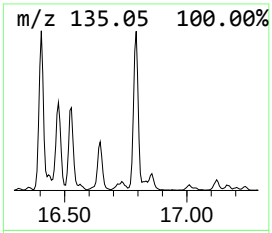
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

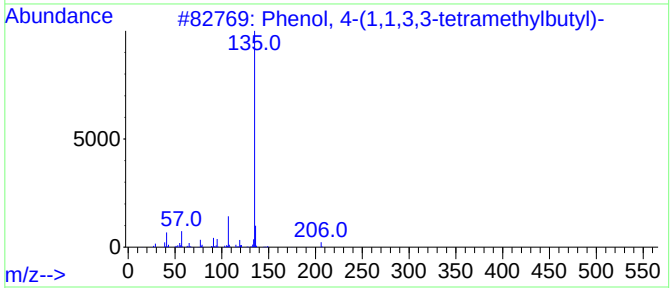
****
Peak Number 5 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 5

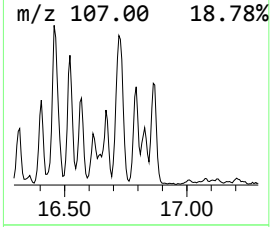
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.792	2.79 ng/ul	421581	Phenanthrene-d10	17.322

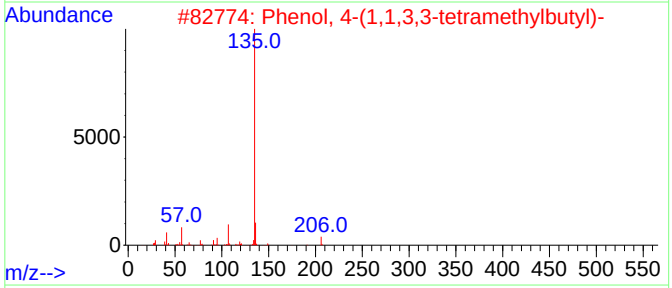
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
3	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	78
4	2-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-37-9	78
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78

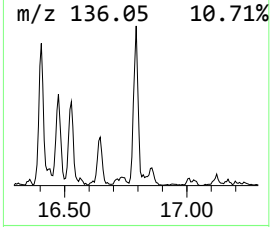


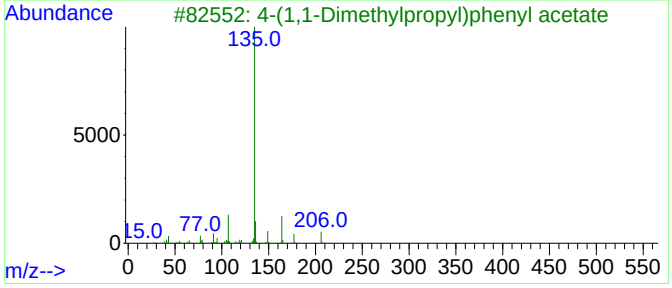


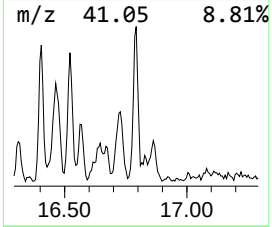


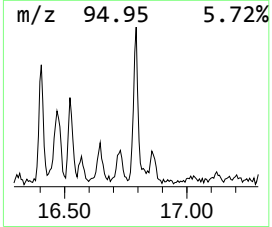












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

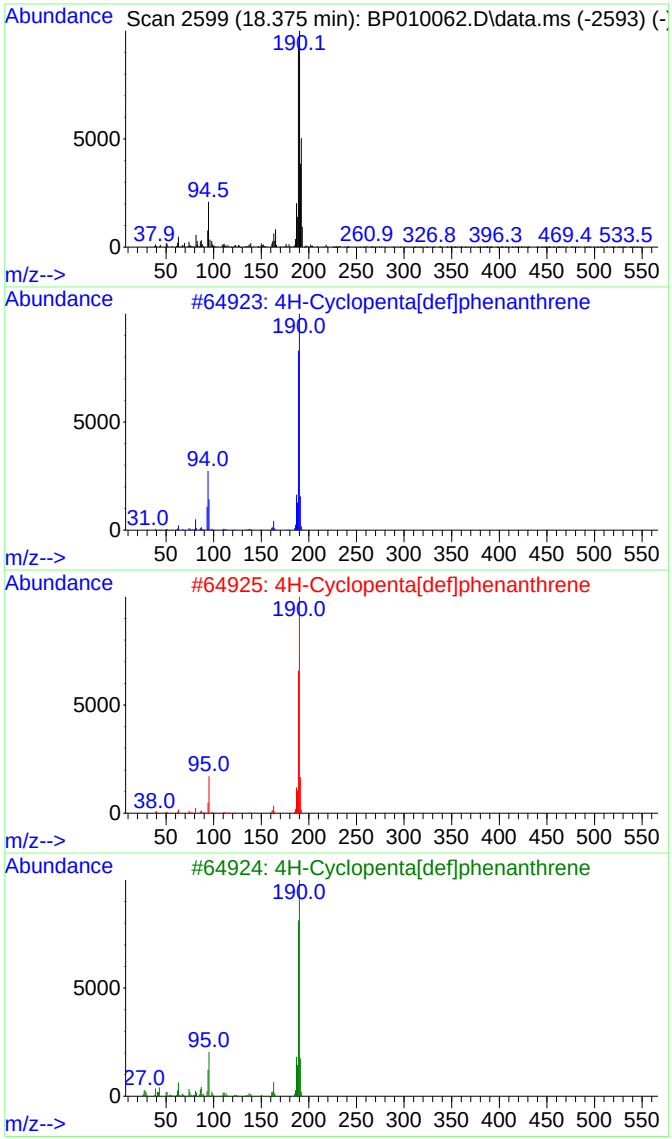
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

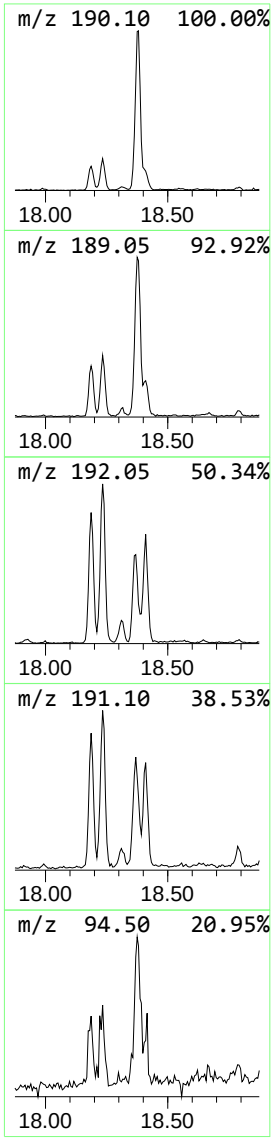
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.375	2.40 ng/ul	361708	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	47
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	22
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	18





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Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

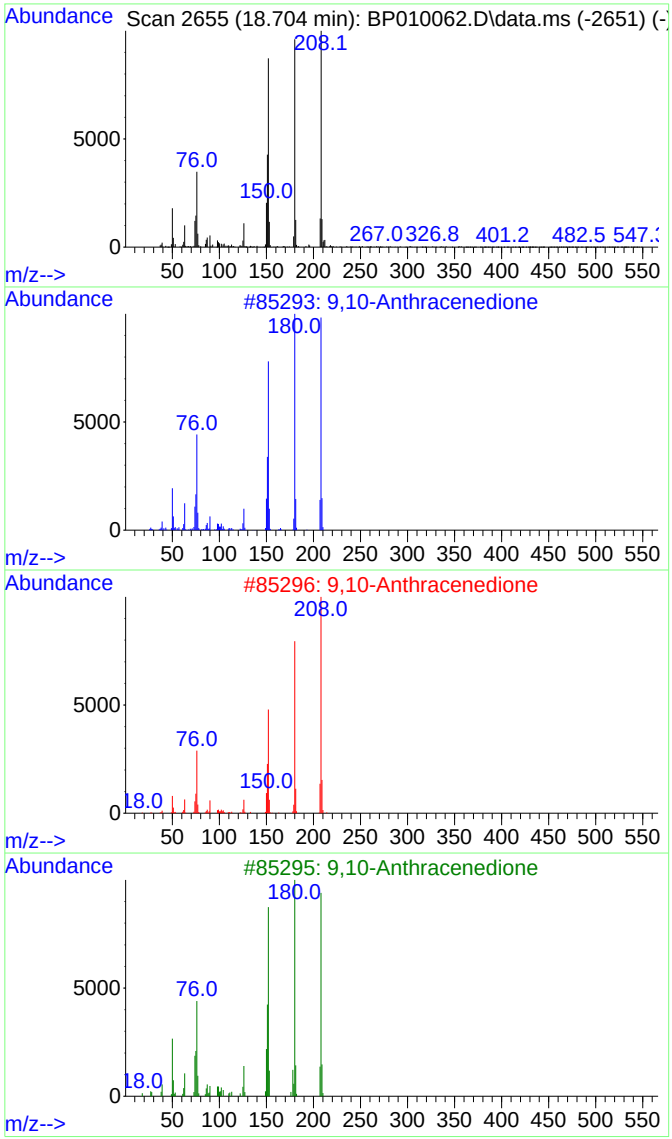
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

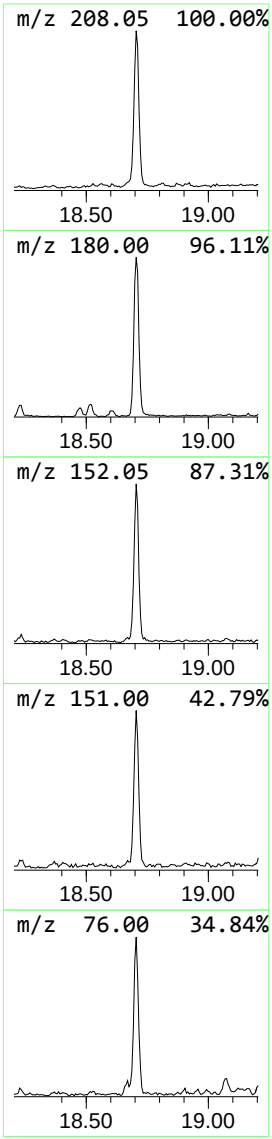
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.704	2.61 ng/ul	394154	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5	9H-Fluoren-9-one	180	C13H8O	000486-25-9	86





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Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

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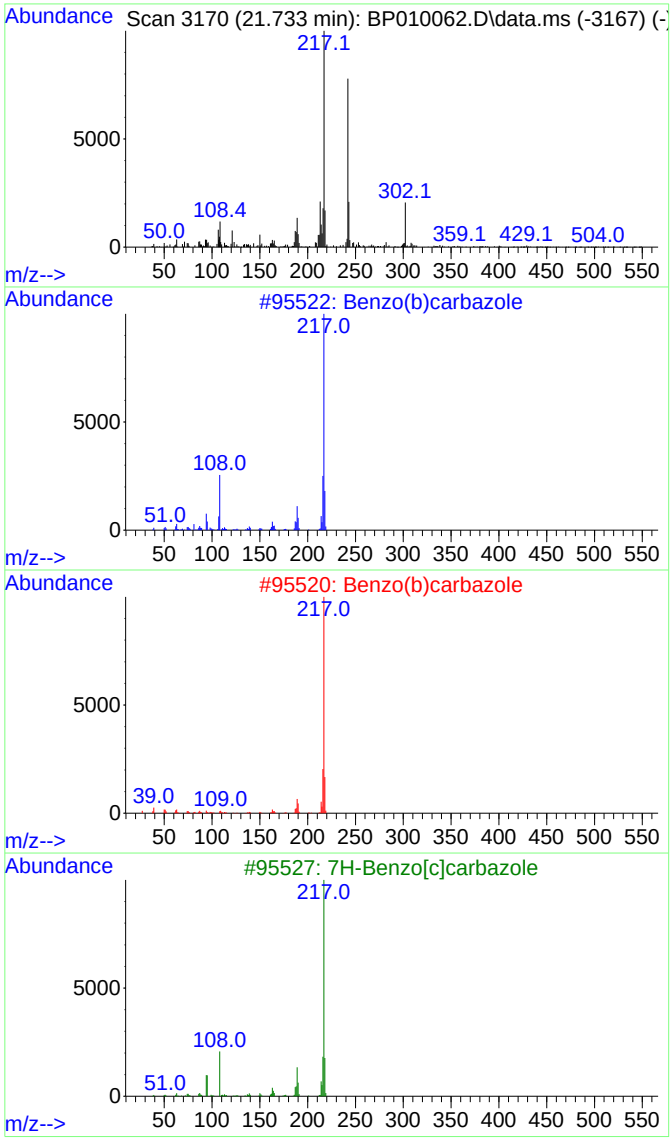
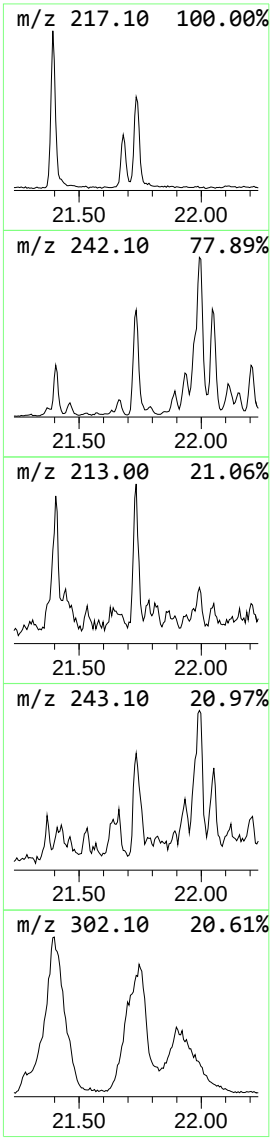
Peak Number 8 Benzo(b)carbazole Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.733	2.48 ng/ul	640815	Chrysene-d12	21.404

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo(b)carbazole	217	C16H11N	000243-28-7	81
2	Benzo(b)carbazole	217	C16H11N	000243-28-7	46
3	7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	46
4	7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	42
5	11H-Indeno(1,2-b)quinoline	217	C16H11N	000243-51-6	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

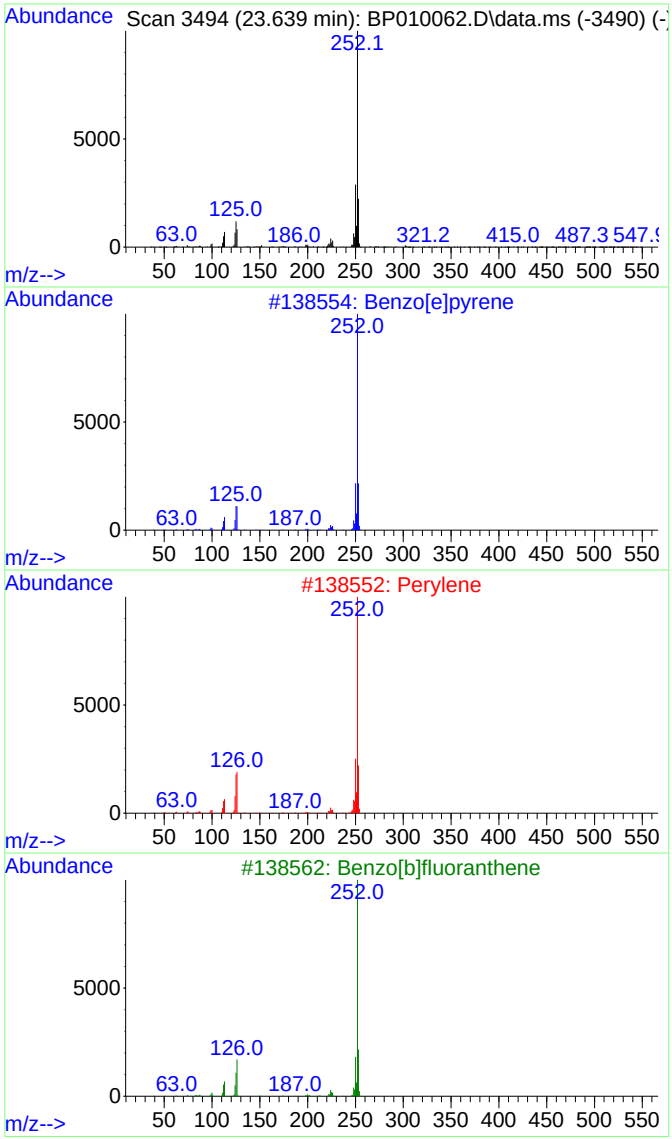
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

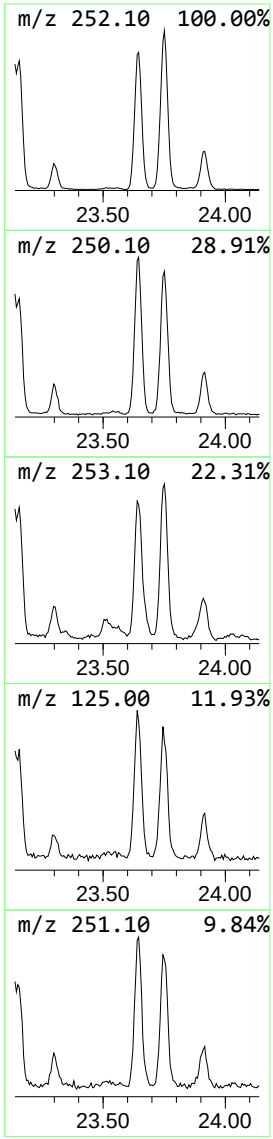
****

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.639	6.56 ng/ul	815450	Perylene-d12	23.857

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2	Perylene	252	C20H12	000198-55-0	94
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4	Perylene	252	C20H12	000198-55-0	93
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010062.D
Acq On : 22 Apr 2022 19:26
Operator : CG/JU
Sample : N2373-08DL 5X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFE0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.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
Phenol, 2-(1,1,...	16.404	2.6	ng/ul	387242	4	17.322	3017130	20.0
4-(7-Methylocty...	16.463	3.8	ng/ul	568396	4	17.322	3017130	20.0
Phenol, p-tert-...	16.522	3.0	ng/ul	444626	4	17.322	3017130	20.0
Acetamide, N-(3...	16.722	3.1	ng/ul	461491	4	17.322	3017130	20.0
Phenol, 4-(1,1,...	16.792	2.8	ng/ul	421581	4	17.322	3017130	20.0
4H-Cyclopenta[d...	18.375	2.4	ng/ul	361708	4	17.322	3017130	20.0
9,10-Anthracene...	18.704	2.6	ng/ul	394154	4	17.322	3017130	20.0
Benzo(b)carbazole	21.733	2.5	ng/ul	640815	5	21.404	5169660	20.0
Benzo[e]pyrene	23.639	6.6	ng/ul	815450	6	23.857	2486180	20.0