LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCYN8

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-BP100724.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP022355.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.223	37	40	47	rVB	17439	21766	0.56%	0.046%
2	3.634	107	110	126	rBV	181298	282001	7.19%	0.593%
3	3.958	160	165	173	rVB4	8087	12748	0.33%	0.027%
4	6.887	652	663	674	rBV	303584	511805	13.05%	1.076%
5	7.040	683	689	698	rVB	222917	358447	9.14%	0.754%
6	7.240	715	723	731	rBV	303871	509621	13.00%	1.072%
7	7.305	731	734	741	rVB2	10825	17078	0.44%	0.036%
8	7.493	758	766	773	rBV	24307	46009	1.17%	0.097%
9	7.699	793	801	809	rBV	441081	712957	18.18%	1.499%
10	8.399	908	920	932	rBV	413313	684083	17.44%	1.439%
11	8.834	988	994	1003	rBV	291696	505779	12.90%	1.064%
12	9.252	1055	1065	1069	rBV3	7245	13415	0.34%	0.028%
13	9.393	1080	1089	1098	rBV2	26121	46639	1.19%	0.098%
14	9.552	1108	1116	1124	rBV	276766	459193	11.71%	0.966%
15	10.093	1196	1208	1220	rBV	434209	756720	19.30%	1.591%
16	10.463	1262	1271	1276	rBV	559434	1017741	25.95%	2.140%
17	10.510	1276	1279	1285	rVV	101167	148453	3.79%	0.312%
18	10.593	1285	1293	1310	rVB	366010	679786	17.34%	1.429%
19	10.999	1355	1362	1368	rBV8	9632	19492	0.50%	0.041%
20	12.034	1535	1538	1545	rVV3	8166	14400	0.37%	0.030%
21	12.116	1545	1552	1560	rVB	117087	199792	5.09%	0.420%
22	12.334	1581	1589	1598	rVB	81381	132941	3.39%	0.280%
23	13.134	1719	1725	1733	rBV	85905	128825	3.29%	0.271%
24	13.334	1753	1759	1769	rVB	25402	51888	1.32%	0.109%
25	13.463	1774	1781	1782	rBV	44205	48283	1.23%	0.102%
26	13.628	1801	1809	1814	rBV2	58101	118764	3.03%	0.250%
27	13.722	1814	1825	1834	rVB	835762	1275124	32.52%	2.681%
28	13.857	1844	1848	1852	rBV	28547	40449	1.03%	0.085%
29	13.893	1852	1854	1862	rVB3	10637	14640	0.37%	0.031%
30	14.004	1865	1873	1891	rBV2	872340	1538772	39.24%	3.236%
31	14.316	1915	1926	1933	rBV	976814	1645535	41.96%	3.460%
32	14.387	1933	1938	1945	rVV	252102	425021	10.84%	0.894%
33	14.446	1945	1948	1955	rVB	33709	52042	1.33%	0.109%
34	14.540	1955	1964	1974	rBV	276334	499071	12.73%	1.049%
35	14.622	1976	1978	1983	rVV4	13621	23619	0.60%	0.050%
36	14.722	1984	1995	2004	rVV	399445	701485	17.89%	1.475%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCYN8

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-BP100724.MA.M
Title : SVOA CALIBRATION

37	14.810	2005	2010	2016	rVB2	57549	86493	2.21%	0.182%
38	14.928	2024	2030	2037	rBV4	21421	50173	1.28%	0.106%
39	14.987	2037	2040	2048	rVB3	17765	31475	0.80%	0.066%
40	15.128	2058	2064	2066	rVV2	17520	29003	0.74%	0.061%
41	15.157	2066	2069	2076	rVB6	15404	28500	0.73%	0.060%
42	15.316	2089	2096	2102	rVV	1242382	1981487	50.53%	4.167%
43	15.381	2102	2107	2112	rVV	347807	566070	14.44%	1.190%
44	15.457	2115	2120	2127	rVV	290381	438914	11.19%	0.923%
45	15.510	2127	2129	2132	rVV	37747	49354	1.26%	0.104%
46	15.545	2132	2135	2140	rVV	68268	102321	2.61%	0.215%
47	15.587	2140	2142	2144	rVV3	14219	15885	0.41%	0.033%
48	15.622	2144	2148	2152	rVV2	28605	49711	1.27%	0.105%
49	15.698	2152	2161	2169	rVV2	472618	889452	22.68%	1.870%
50	15.834	2177	2184	2191	rVV	139453	273741	6.98%	0.576%
51	15.910	2194	2197	2204	rVB2	31134	51860	1.32%	0.109%
52	16.069	2212	2224	2231	rBV7	23785	80151	2.04%	0.169%
53	16.269	2253	2258	2263	rBV	34011	47137	1.20%	0.099%
54	16.410	2269	2282	2287	rBV3	53212	125624	3.20%	0.264%
55	16.457	2287	2290	2294	rVB	19489	26601	0.68%	0.056%
56	16.510	2294	2299	2301	rBV6	10199	13433	0.34%	0.028%
57	16.557	2301	2307	2312	rVB3	17874	33112	0.84%	0.070%
58	16.640	2313	2321	2324	rBV2	34337	65230	1.66%	0.137%
59	16.675	2324	2327	2331	rVB3	29671	33110	0.84%	0.070%
60	16.763	2338	2342	2349	rVB3	31334	56838	1.45%	0.120%
61	16.863	2351	2359	2363	rVV5	40824	107773	2.75%	0.227%
62	16.922	2363	2369	2388	rVB	187953	343660	8.76%	0.723%
63	17.116	2395	2402	2405	rBV	1381187	2118433	54.02%	4.455%
64	17.163	2405	2410	2414	rVB	2621337	3921371	100.00%	8.246%
65	17.216	2414	2419	2425	rBV	1351091	2139022	54.55%	4.498%
66	17.310	2431	2435	2440	rVB4	25092	48441	1.24%	0.102%
67	17.545	2469	2475	2481	rBV	115623	199457	5.09%	0.419%
68	17.651	2488	2493	2499	rVB3	40096	64039	1.63%	0.135%
69	17.710	2499	2503	2507	rBV2	25605	37476	0.96%	0.079%
70	17.863	2525	2529	2534	rVB3	21545	33767	0.86%	0.071%
71	18.022	2547	2556	2559	rBV	202520	317856	8.11%	0.668%
72	18.063	2559	2563	2572	rVV2	518974	862147	21.99%	1.813%
73	18.151	2574	2578	2584	rVV	60210	104620	2.67%	0.220%
74	18.228	2584	2591	2595	rVV3	189982	396140	10.10%	0.833%
75	18.263	2595	2597	2603	rVB	93595	123748	3.16%	0.260%
76	18.551	2640	2646	2650	rBV	123707	189356	4.83%	0.398%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCYN8

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-BP100724.MA.M
Title : SVOA CALIBRATION

77	18.592	2650	2653	2657	rVB3	24561	35122	0.90%	0.074%
78	18.792	2684	2687	2691	rVB3	31892	43555	1.11%	0.092%
79	18.851	2694	2697	2700	rVV	19932	23080	0.59%	0.049%
80	18.892	2701	2704	2707	rVV2	18119	23117	0.59%	0.049%
81	18.981	2715	2719	2723	rBV	38454	55482	1.41%	0.117%
82	19.028	2724	2727	2728	rBV3	23014	25771	0.66%	0.054%
83	19.122	2738	2743	2752	rVB7	35480	93566	2.39%	0.197%
84	19.251	2756	2765	2773	rBV	1724038	2492867	63.57%	5.242%
85	19.386	2785	2788	2794	rVB2	68362	102786	2.62%	0.216%
86	19.498	2804	2807	2812	rVB2	37077	48287	1.23%	0.102%
87	19.598	2818	2824	2828	rBV	2120480	3470893	88.51%	7.299%
88	19.634	2828	2830	2837	rVV	1044207	1157185	29.51%	2.433%
89	19.710	2839	2843	2848	rVB2	48723	75269	1.92%	0.158%
90	19.810	2857	2860	2865	rVB	66517	92578	2.36%	0.195%
91	19.963	2882	2886	2888	rBV2	49249	77169	1.97%	0.162%
92	20.145	2913	2917	2922	rVB	119542	168921	4.31%	0.355%
93	20.245	2931	2934	2939	rVB	103748	158016	4.03%	0.332%
94	21.128	3081	3084	3088	rBV	55465	82287	2.10%	0.173%
95	21.169	3089	3091	3095	rVB2	61650	56861	1.45%	0.120%
96	21.228	3096	3101	3106	rVV2	285389	459045	11.71%	0.965%
97	21.563	3150	3158	3163	rBV2	1772272	3070923	78.31%	6.458%
98	21.604	3163	3165	3175	rVB	232541	328584	8.38%	0.691%
99	24.639	3672	3681	3689	rBV	1088975	2778432	70.85%	5.843%
100	24.857	3710	3718	3731	rVB	986957	2787794	71.09%	5.862%

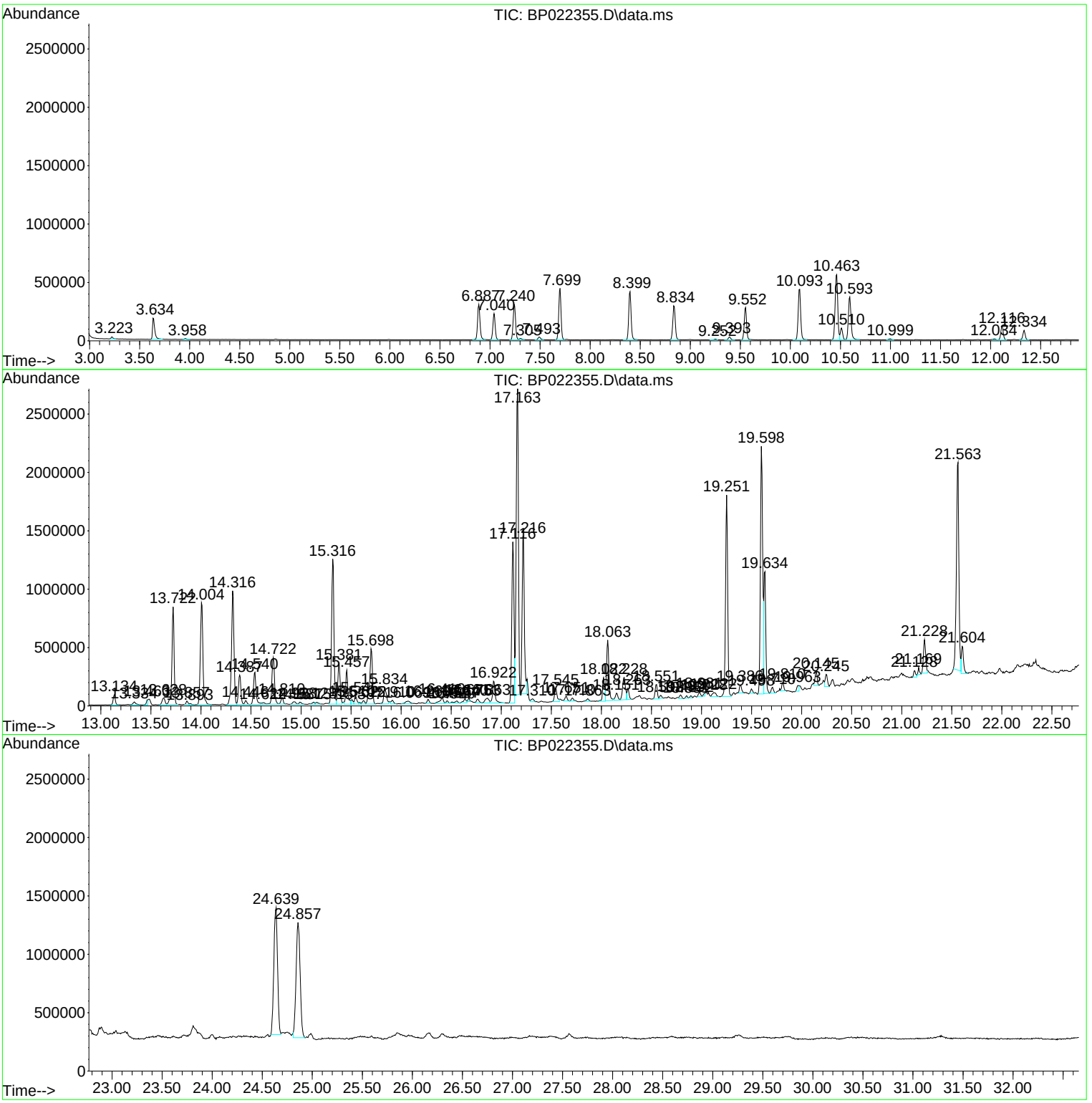
Sum of corrected areas: 47554955

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

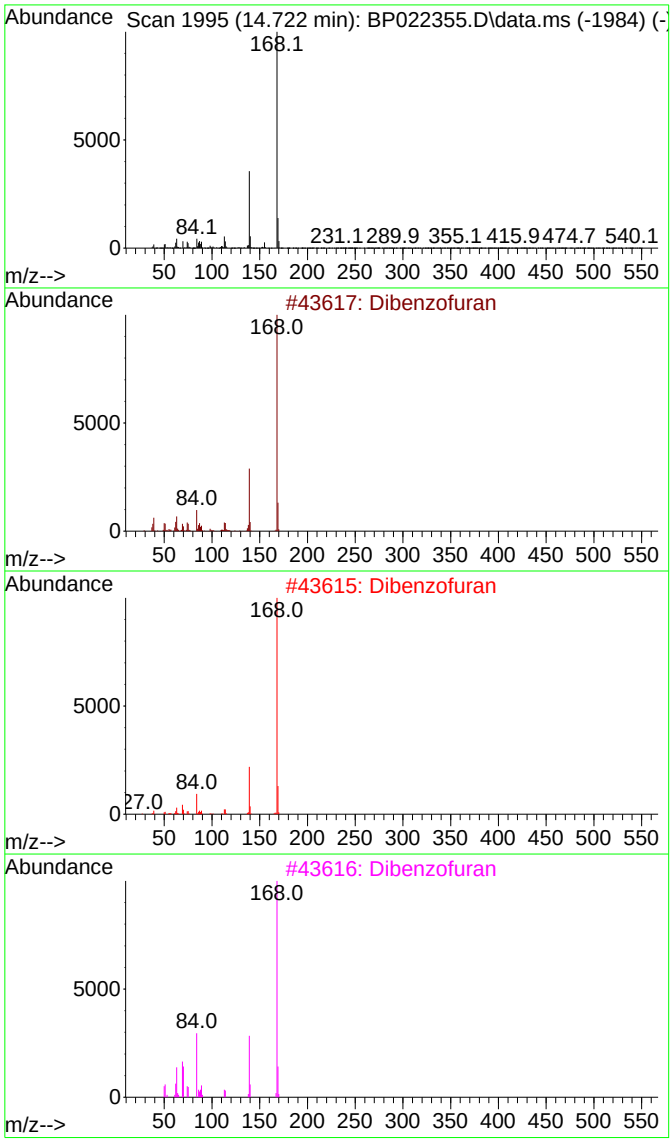
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

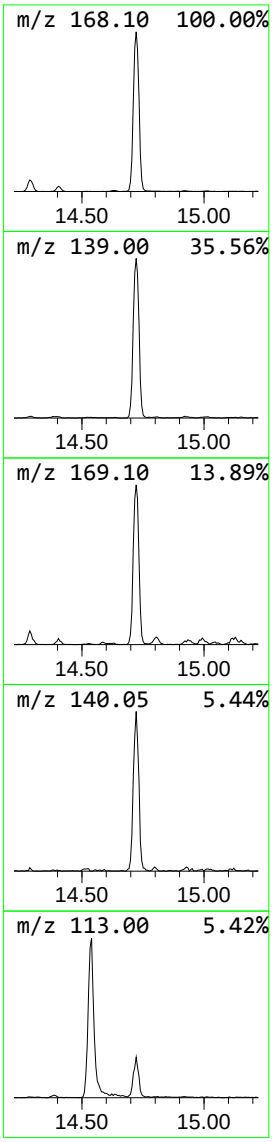
****

Peak Number 1 Dibenzo furan Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.722	8.53 ng/ul	701485	Acenaphthene-d10	14.316

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	87
2	Dibenzofuran	168	C12H8O	000132-64-9	80
3	Dibenzofuran	168	C12H8O	000132-64-9	78
4	Naphtho[2,1-d]imidazole	168	C11H8N2	000233-53-4	59
5	1H-Pyrido[2,3-b]indole	168	C11H8N2	000244-76-8	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

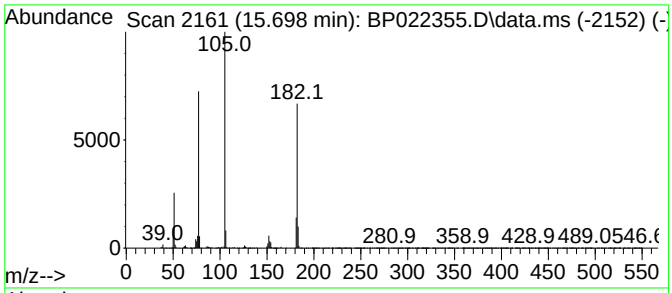
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

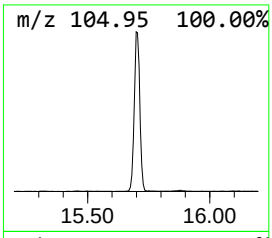
****

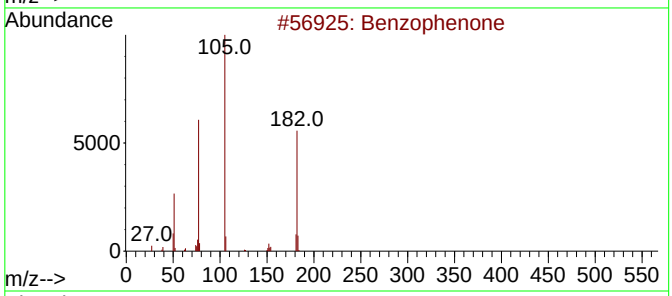
Peak Number 2 Benzophenone Concentration Rank 1

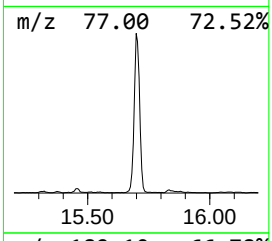
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.698	10.81 ng/ul	889452	Acenaphthene-d10	14.316

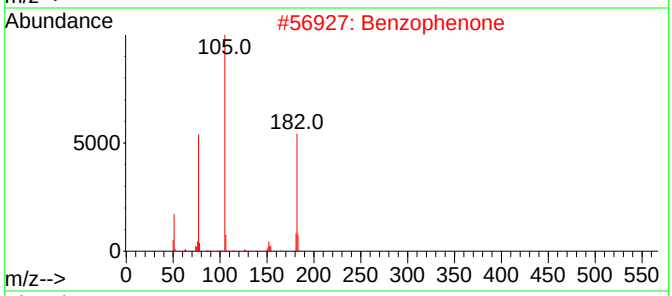
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	87
5	1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64

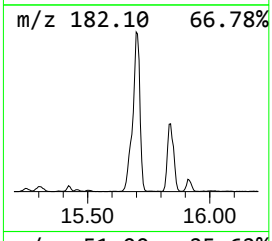


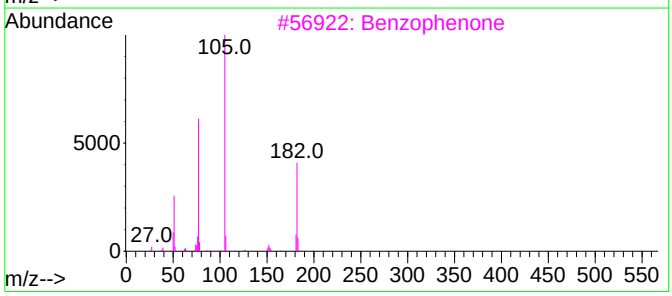


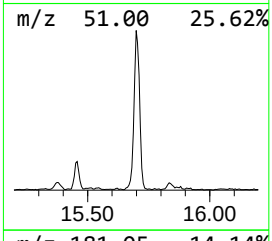


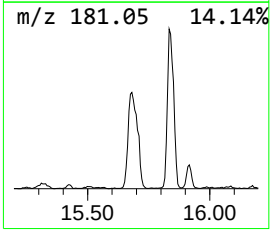












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

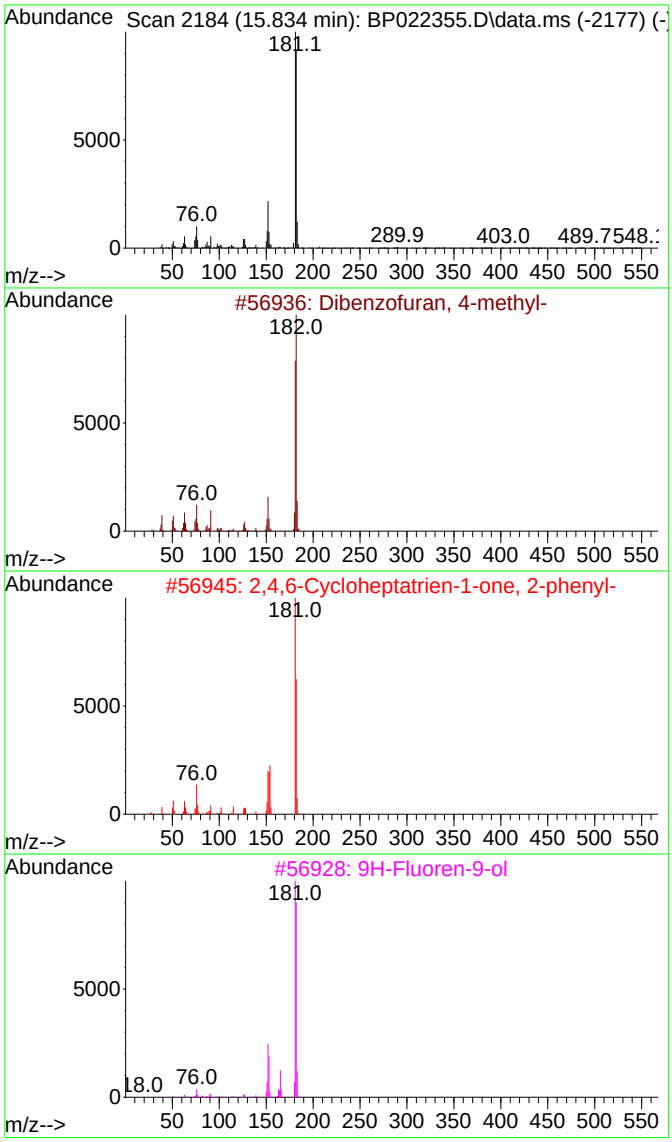
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

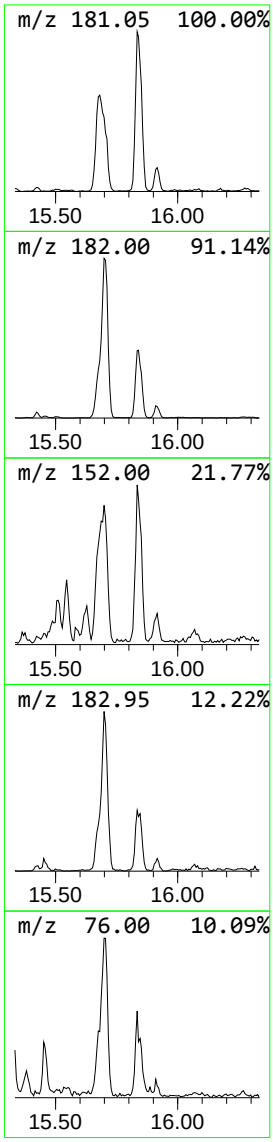
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Dibenzofuran, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.834	2.58 ng/ul	273741	Phenanthrene-d10	17.116

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	90
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	90
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
5	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

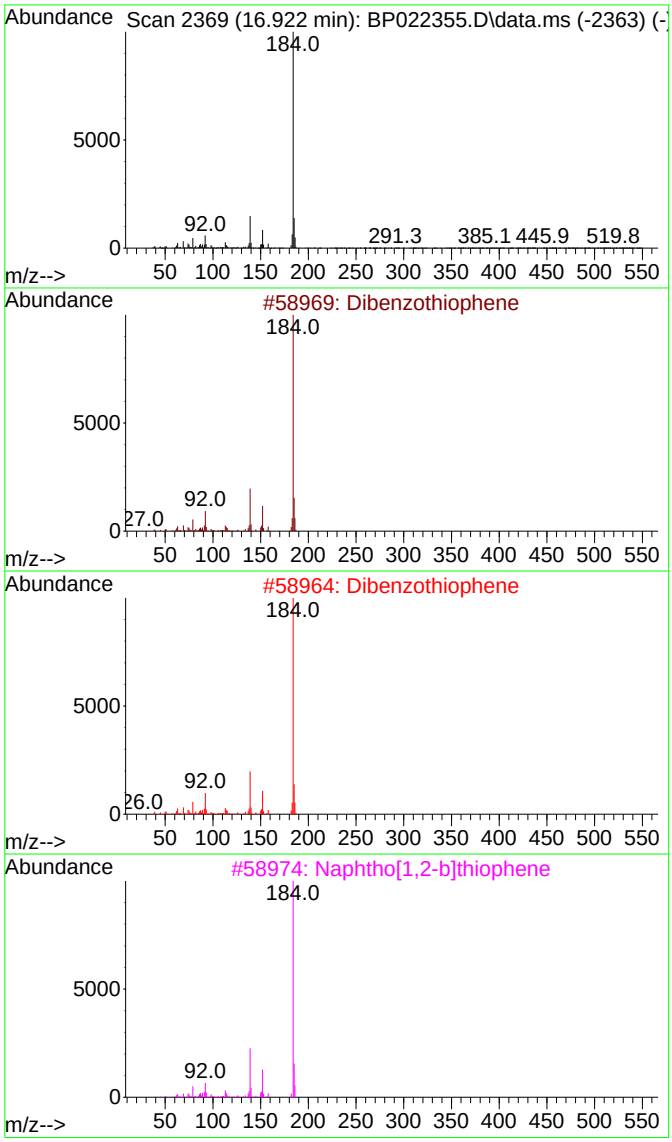
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

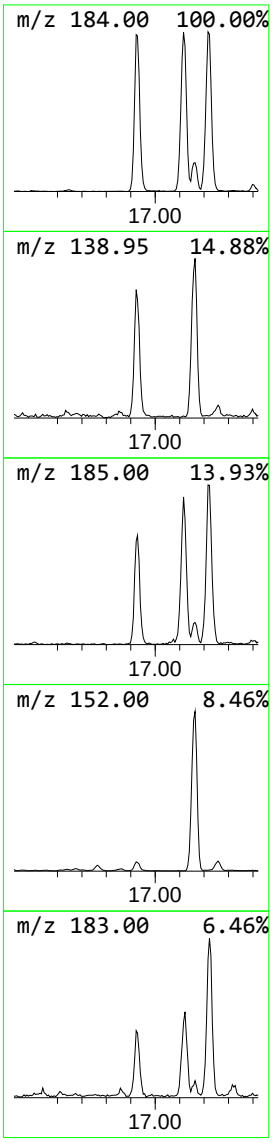
****

Peak Number 4 Dibenzothiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.922	3.24 ng/ul	343660	Phenanthrene-d10	17.116

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[1,2-b]thiophene	184	C12H8S	000234-41-3	95
4	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
5	Dibenzothiophene	184	C12H8S	000132-65-0	95





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Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

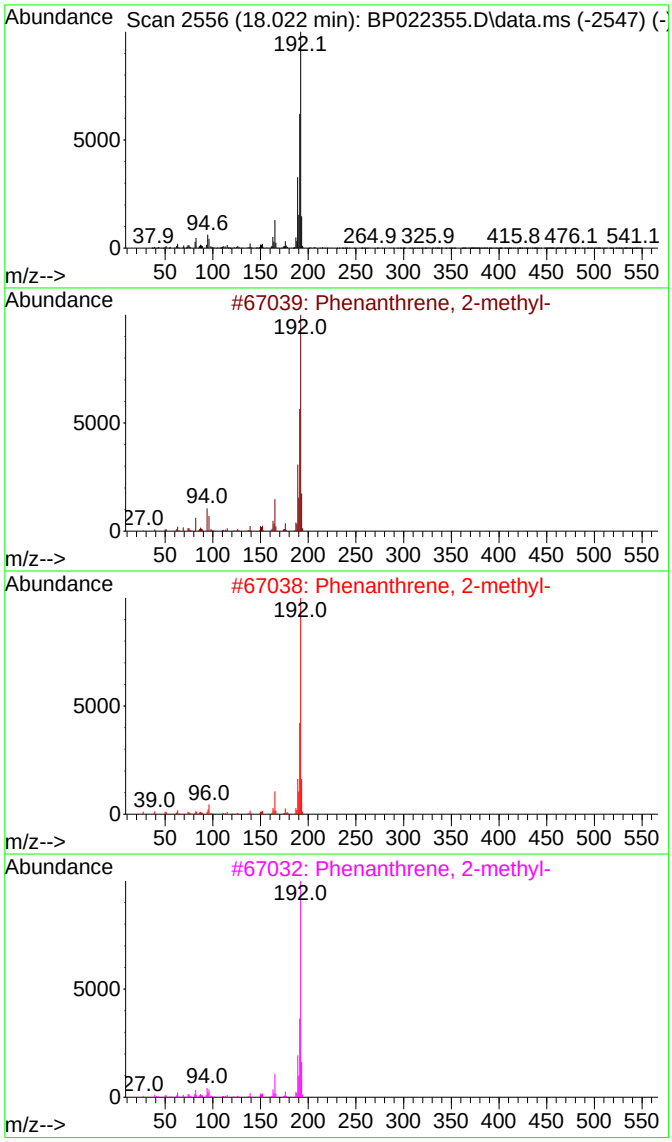
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

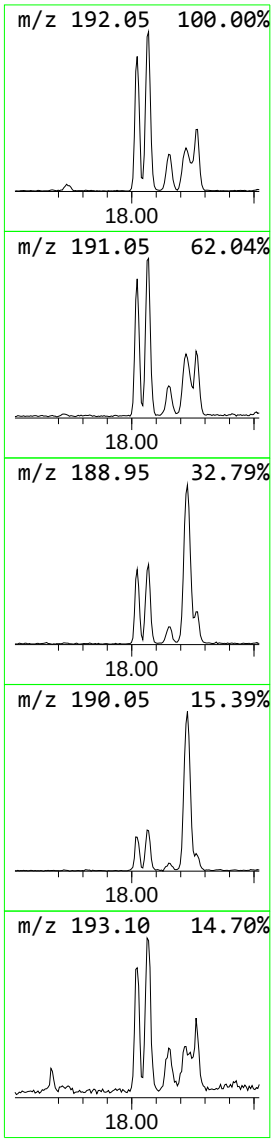
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	3.00 ng/ul	317856	Phenanthrene-d10	17.116

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





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Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

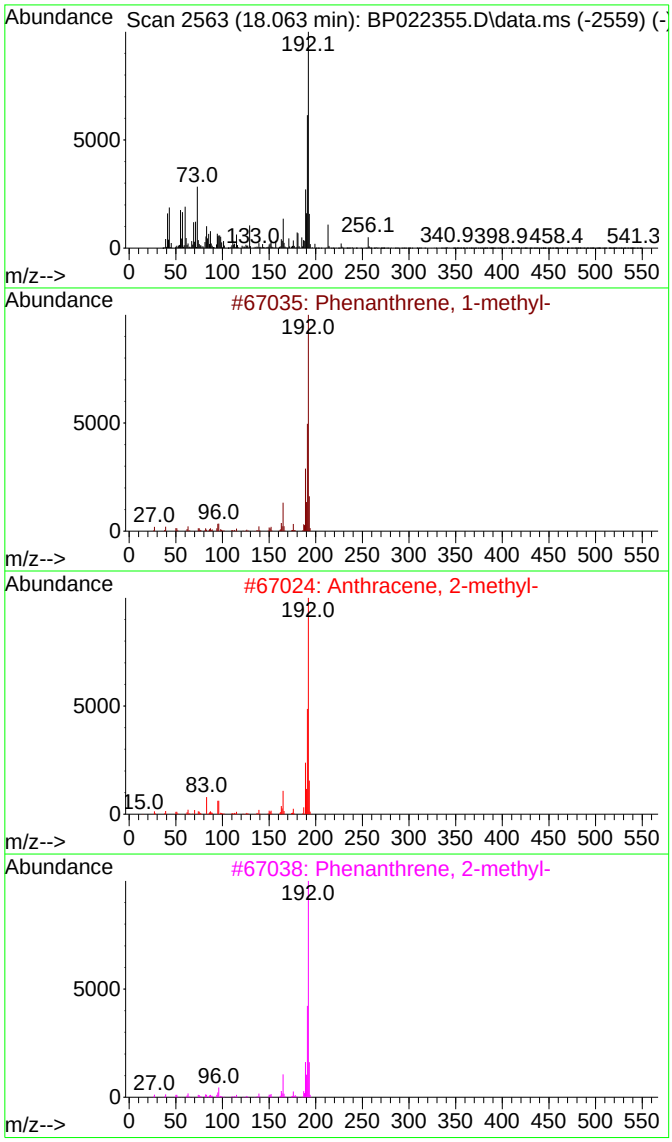
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

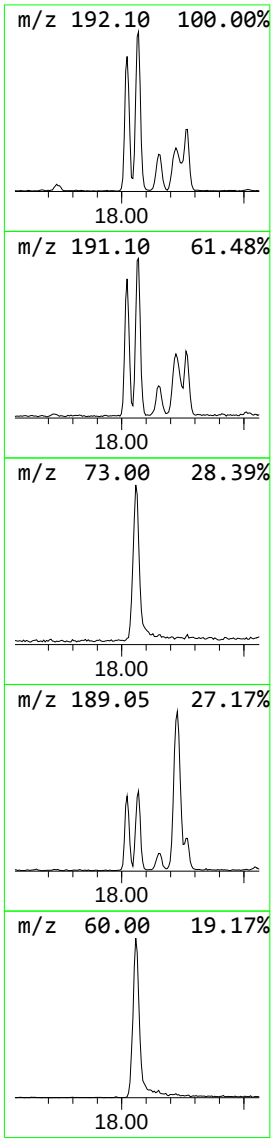
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.063	8.14 ng/ul	862147	Phenanthrene-d10	17.116

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	92





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Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

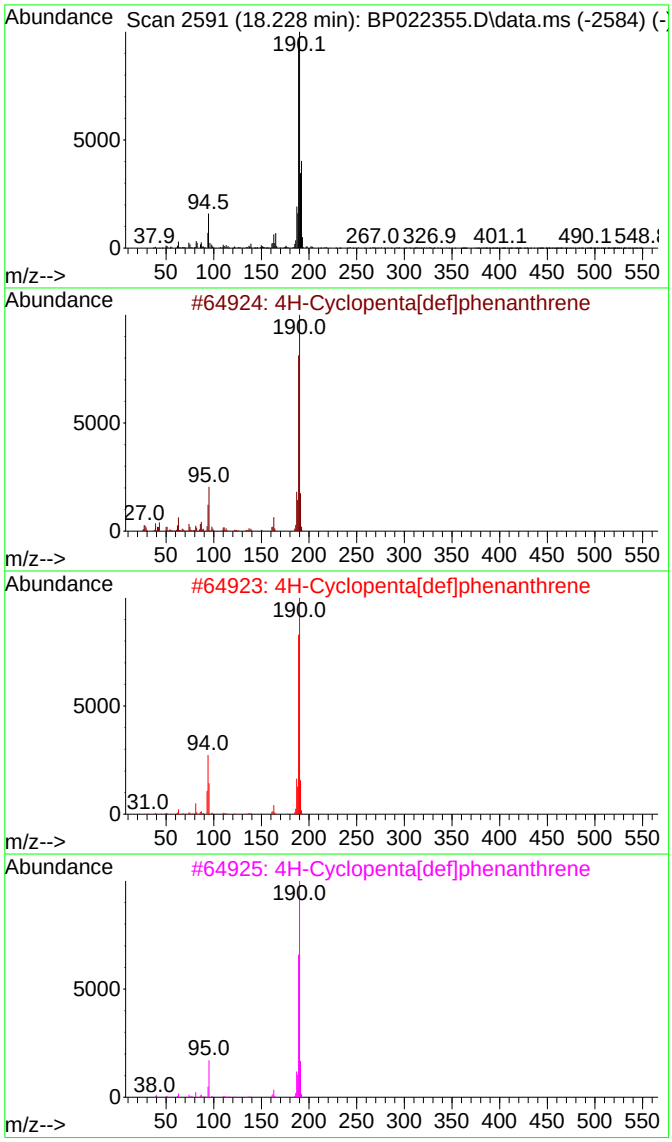
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

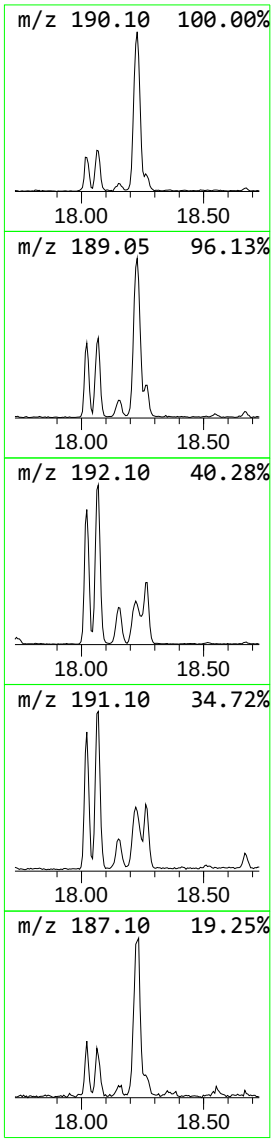
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	3.74 ng/ul	396140	Phenanthrene-d10	17.116

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	70
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
5	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

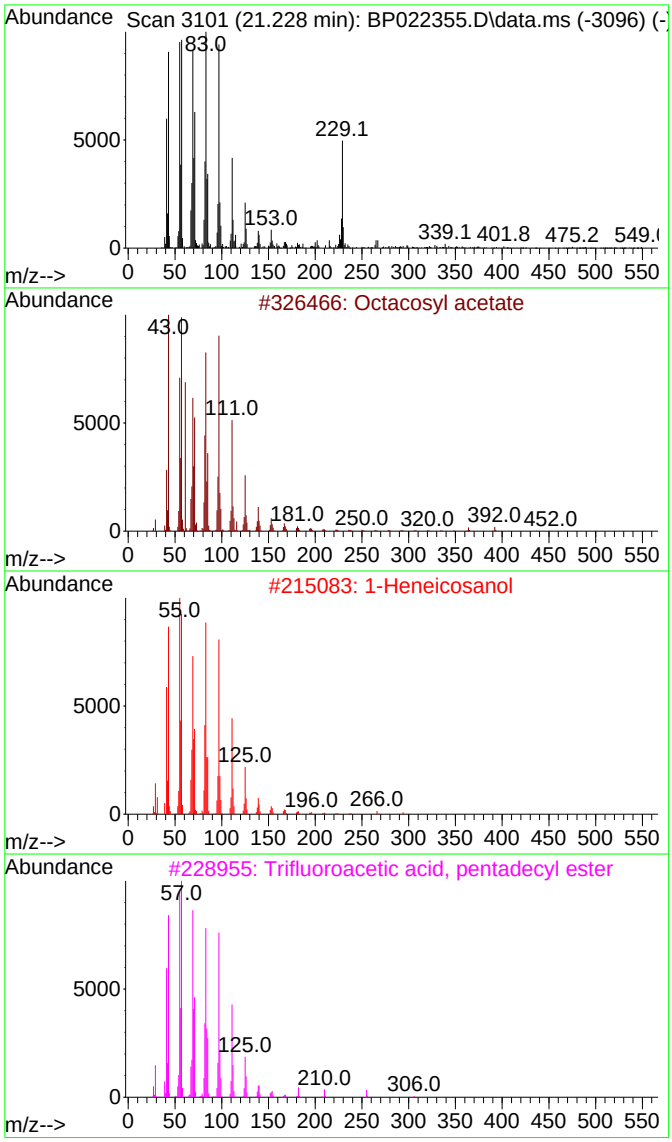
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

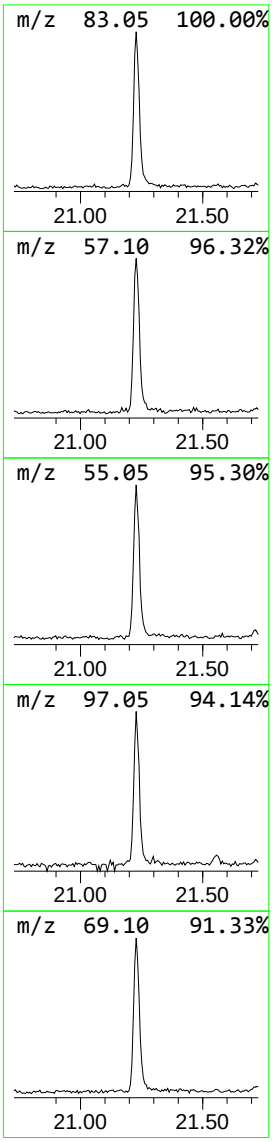
****

Peak Number 8 Octacosyl acetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.227	2.99 ng/ul	459045	Chrysene-d12	21.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosyl acetate	452	C30H60O2	018206-97-8	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
Data File : BP022355.D
Acq On : 14 Oct 2024 14:35
Operator : RC/JU
Sample : P4264-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYN8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Dibenzo furan	14.722	8.5	ng/ul	701485	3	14.316	1645540	20.0
Benzophenone	15.698	10.8	ng/ul	889452	3	14.316	1645540	20.0
Dibenzofuran, 4...	15.834	2.6	ng/ul	273741	4	17.116	2118430	20.0
Dibenzothiophene	16.922	3.2	ng/ul	343660	4	17.116	2118430	20.0
Phenanthrene, 2...	18.022	3.0	ng/ul	317856	4	17.116	2118430	20.0
Phenanthrene, 1...	18.063	8.1	ng/ul	862147	4	17.116	2118430	20.0
4H-Cyclopenta[d...	18.228	3.7	ng/ul	396140	4	17.116	2118430	20.0
Octacosyl acetate	21.227	3.0	ng/ul	459045	5	21.563	3070920	20.0