LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A43N1

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-BP050124.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP020242.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	18	22	31	rVB4	17800	27272	0.78%	0.091%
2	3.481	64	67	83	rBV	54508	99312	2.84%	0.332%
3	3.622	87	91	103	rVV	100313	185731	5.30%	0.620%
4	3.711	103	106	114	rVB3	18069	28148	0.80%	0.094%
5	3.905	136	139	147	rVB	7134	12924	0.37%	0.043%
6	4.311	204	208	213	rVB3	2954	4520	0.13%	0.015%
7	4.405	220	224	229	rBV7	2653	5124	0.15%	0.017%
8	4.452	229	232	235	rVB5	3577	3966	0.11%	0.013%
9	4.675	266	270	275	rBV5	4441	7363	0.21%	0.025%
10	4.793	284	290	296	rBV2	9418	17535	0.50%	0.059%
11	5.134	342	348	353	rBV2	5950	9273	0.26%	0.031%
12	5.658	431	437	439	rBV7	3408	5132	0.15%	0.017%
13	5.910	476	480	486	rVB6	5486	9062	0.26%	0.030%
14	6.669	605	609	613	rVV7	2321	4445	0.13%	0.015%
15	6.722	614	618	623	rVB6	3489	5856	0.17%	0.020%
16	6.805	623	632	650	rBV	245135	511572	14.61%	1.708%
17	6.975	654	661	673	rVV	248684	428187	12.23%	1.430%
18	7.158	682	692	700	rBV	298601	538453	15.38%	1.798%
19	7.263	707	710	723	rBV2	12741	27578	0.79%	0.092%
20	7.616	764	770	784	rBV	335480	587452	16.78%	1.961%
21	7.728	784	789	794	rVB8	4700	8571	0.24%	0.029%
22	8.105	848	853	857	rBV7	4049	7595	0.22%	0.025%
23	8.240	870	876	882	rBV8	3357	7164	0.20%	0.024%
24	8.334	885	892	907	rBV	165248	358981	10.25%	1.199%
25	8.452	907	912	921	rVB	56553	100825	2.88%	0.337%
26	8.781	960	968	982	rBV	341262	620572	17.72%	2.072%
27	9.128	1022	1027	1032	rBV5	3530	6359	0.18%	0.021%
28	9.357	1060	1066	1078	rBV3	21186	49635	1.42%	0.166%
29	9.493	1081	1089	1105	rBV	295964	543317	15.52%	1.814%
30	9.740	1126	1131	1134	rBV7	4096	9193	0.26%	0.031%
31	9.869	1148	1153	1159	rVB8	4418	8662	0.25%	0.029%
32	10.028	1173	1180	1207	rBV	295267	720163	20.57%	2.404%
33	10.387	1233	1241	1256	rBV	489609	861437	24.60%	2.876%
34	10.551	1262	1269	1285	rBV	235814	543167	15.51%	1.813%
35	10.987	1338	1343	1345	rBV6	7334	14198	0.41%	0.047%
36	11.169	1367	1374	1389	rBV2	42396	137778	3.93%	0.460%

LSC Area Percent Report

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Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

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-BP050124.MA.M
Title : SVOA CALIBRATION

37	11.945	1499	1506	1511	rBV9	2029	4911	0.14%	0.016%
38	12.004	1511	1516	1521	rVB3	5000	10338	0.30%	0.035%
39	12.869	1659	1663	1665	rBV4	3656	4917	0.14%	0.016%
40	13.092	1696	1701	1705	rBV8	2468	4624	0.13%	0.015%
41	13.169	1708	1714	1722	rVV	42421	75129	2.15%	0.251%
42	13.469	1756	1765	1769	rBV9	3383	7404	0.21%	0.025%
43	13.604	1783	1788	1792	rBV8	2570	4565	0.13%	0.015%
44	13.716	1792	1807	1821	rBV	713016	1200845	34.30%	4.009%
45	13.981	1844	1852	1868	rVV	801940	1399127	39.96%	4.671%
46	14.198	1886	1889	1894	rVB6	2586	4188	0.12%	0.014%
47	14.292	1896	1905	1915	rBV	675912	1189689	33.98%	3.972%
48	14.392	1920	1922	1928	rVB7	2969	4761	0.14%	0.016%
49	14.516	1934	1943	1944	rBV	211330	269170	7.69%	0.899%
50	14.534	1944	1946	1949	rVV	236196	343667	9.82%	1.147%
51	14.569	1949	1952	1977	rVV	177403	525218	15.00%	1.754%
52	14.745	1979	1982	1988	rVB8	9636	19762	0.56%	0.066%
53	14.963	2014	2019	2028	rVB2	33849	56402	1.61%	0.188%
54	15.122	2041	2046	2051	rBV4	5547	9305	0.27%	0.031%
55	15.186	2051	2057	2062	rVB10	3823	8403	0.24%	0.028%
56	15.322	2071	2080	2097	rBV	1045355	1731884	49.46%	5.782%
57	15.475	2100	2106	2117	rBV	314530	578875	16.53%	1.933%
58	15.686	2137	2142	2146	rBV8	3806	8674	0.25%	0.029%
59	15.928	2176	2183	2189	rBV8	12089	30656	0.88%	0.102%
60	16.075	2204	2208	2213	rBV7	5656	11174	0.32%	0.037%
61	16.134	2214	2218	2223	rVB6	3411	5928	0.17%	0.020%
62	16.416	2258	2266	2267	rBV8	4717	8995	0.26%	0.030%
63	16.469	2272	2275	2281	rVB8	3267	5226	0.15%	0.017%
64	16.539	2282	2287	2294	rBV4	14730	28814	0.82%	0.096%
65	16.681	2305	2311	2316	rVB5	13653	24334	0.69%	0.081%
66	16.775	2320	2327	2343	rBV	2161517	3501363	100.00%	11.690%
67	16.904	2347	2349	2353	rVB5	6664	7121	0.20%	0.024%
68	17.063	2371	2376	2381	rVB7	6591	13640	0.39%	0.046%
69	17.133	2381	2388	2394	rBV	922209	1476431	42.17%	4.929%
70	17.181	2394	2396	2401	rVV2	91050	128584	3.67%	0.429%
71	17.245	2401	2407	2419	rVV	1098650	1811263	51.73%	6.047%
72	17.345	2419	2424	2429	rVV9	11178	18303	0.52%	0.061%
73	17.457	2439	2443	2448	rVB7	6129	10456	0.30%	0.035%
74	17.610	2458	2469	2475	rBV7	17156	49081	1.40%	0.164%
75	17.680	2477	2481	2485	rBV6	8800	16648	0.48%	0.056%
76	17.780	2493	2498	2507	rVB	176080	250106	7.14%	0.835%

LSC Area Percent Report

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Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A43N1

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-BP050124.MA.M
Title : SVOA CALIBRATION

77	17.863	2507	2512	2525	rVB	50563	71505	2.04%	0.239%
78	17.969	2525	2530	2540	rBV	14443	38469	1.10%	0.128%
79	18.057	2540	2545	2547	rBV2	13536	20887	0.60%	0.070%
80	18.104	2547	2553	2567	rBV	433281	678890	19.39%	2.267%
81	18.269	2575	2581	2584	rBV5	21642	37453	1.07%	0.125%
82	18.298	2584	2586	2589	rVB4	7018	8189	0.23%	0.027%
83	18.410	2602	2605	2608	rBV2	14565	17775	0.51%	0.059%
84	18.539	2624	2627	2631	rBV6	7377	11527	0.33%	0.038%
85	18.598	2633	2637	2642	rVV4	11803	22243	0.64%	0.074%
86	18.698	2651	2654	2658	rVB3	11651	15073	0.43%	0.050%
87	18.804	2669	2672	2676	rBV6	9872	15506	0.44%	0.052%
88	18.933	2691	2694	2699	rVB5	13193	15982	0.46%	0.053%
89	19.022	2705	2709	2714	rBV4	18261	26668	0.76%	0.089%
90	19.075	2715	2718	2720	rBV4	15628	19505	0.56%	0.065%
91	19.104	2720	2723	2727	rVB3	36394	44059	1.26%	0.147%
92	19.186	2732	2737	2741	rVV7	27001	48678	1.39%	0.163%
93	19.292	2750	2755	2766	rVB	174405	281156	8.03%	0.939%
94	19.451	2777	2782	2790	rBV2	190653	299342	8.55%	0.999%
95	19.645	2809	2815	2827	rVB	1497940	2433220	69.49%	8.124%
96	20.780	3005	3008	3012	rVB3	26742	33902	0.97%	0.113%
97	21.333	3097	3102	3109	rBV3	130779	258100	7.37%	0.862%
98	21.633	3146	3153	3160	rBV2	760998	1437303	41.05%	4.799%
99	24.774	3678	3687	3696	rBV	630374	1577078	45.04%	5.265%
100	24.992	3717	3724	3736	rVB3	387670	1163450	33.23%	3.884%

Sum of corrected areas: 29952463

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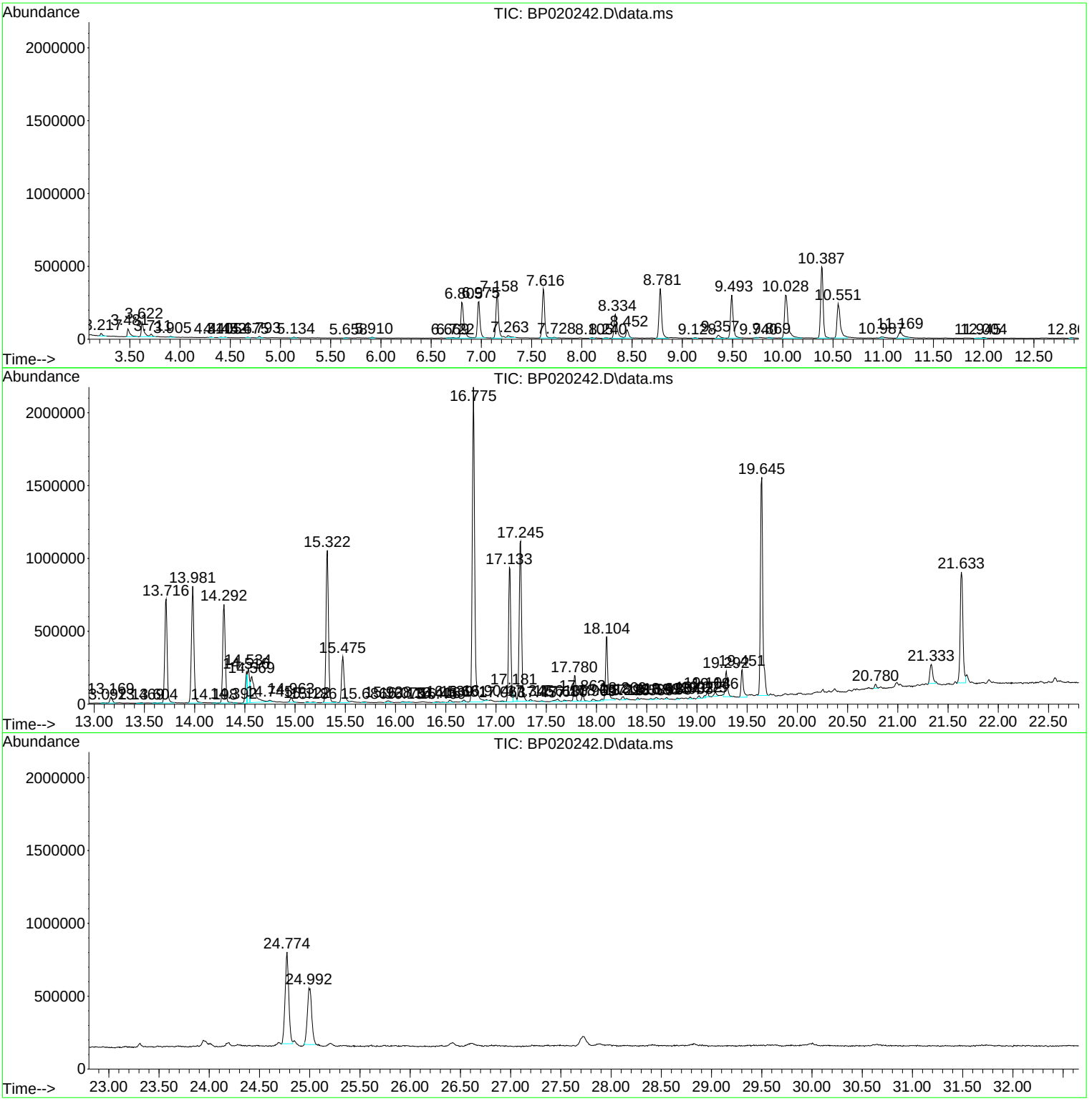
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.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\BP050824\
Data File : BP020242.D
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Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

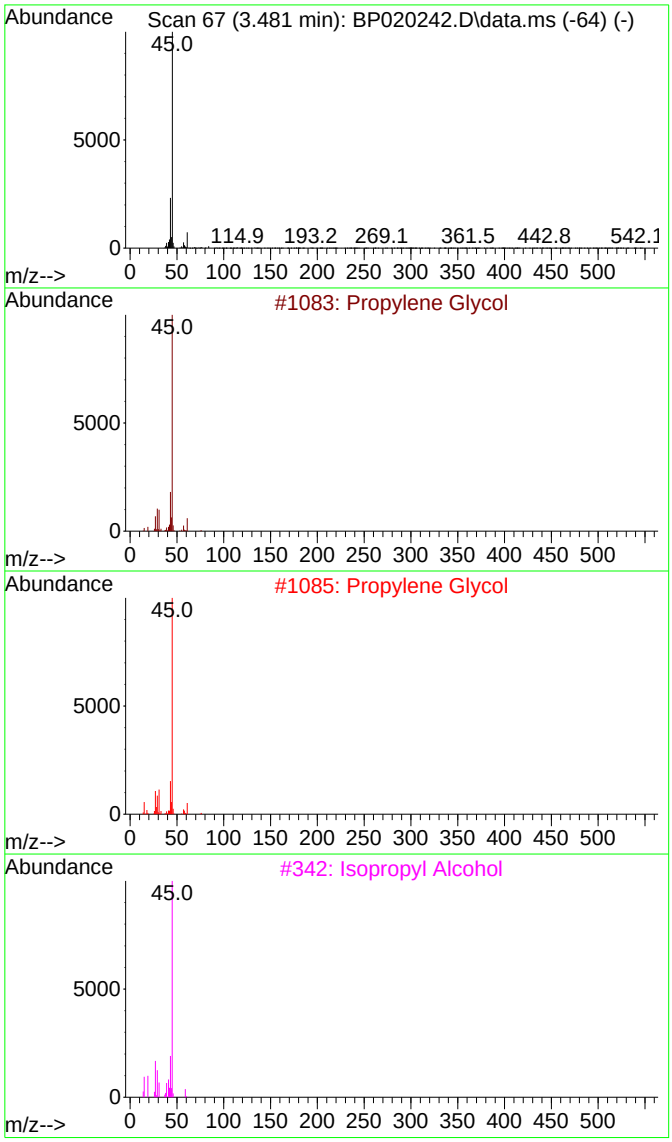
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

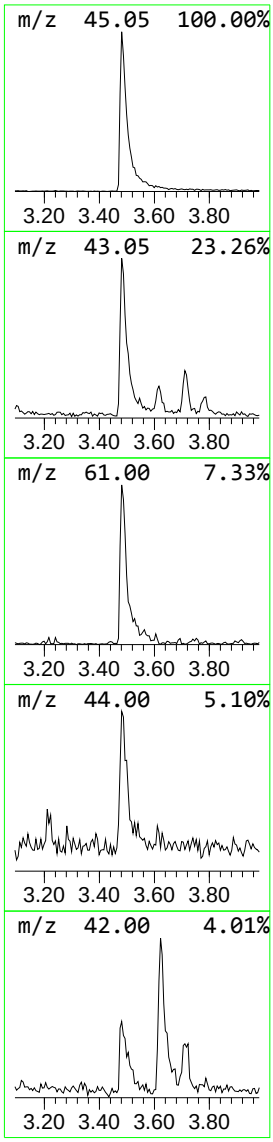
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propylene Glycol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.481	3.38 ng/ul	99312	1,4-Dichlorobenzene-d4	7.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	78
2	Propylene Glycol	76	C3H8O2	000057-55-6	72
3	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

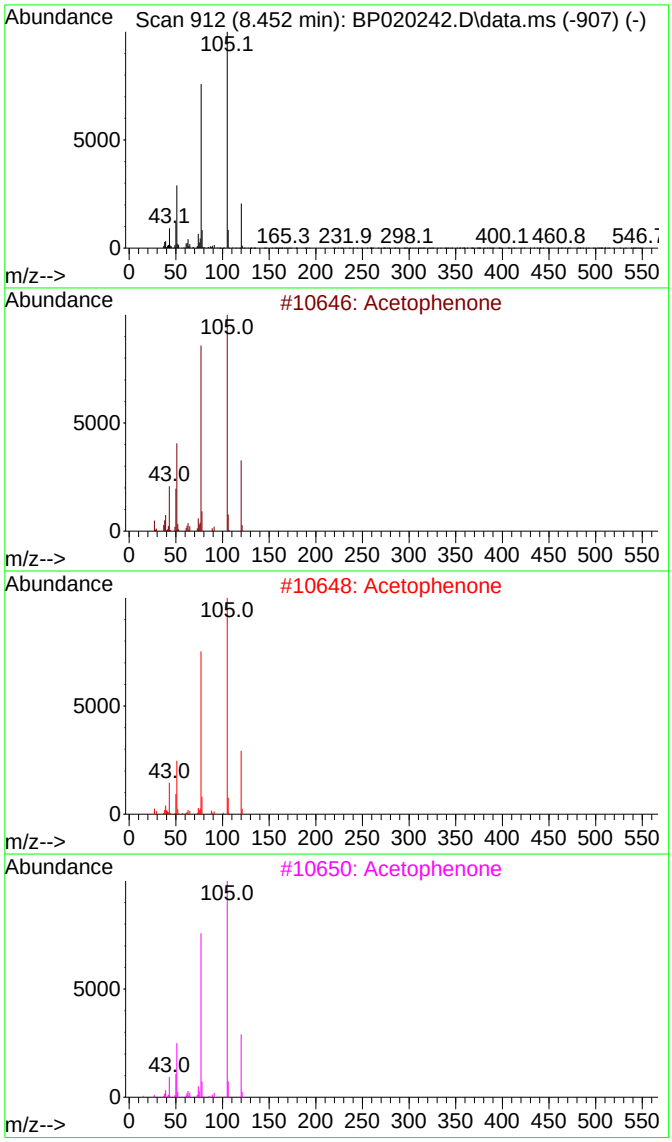
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

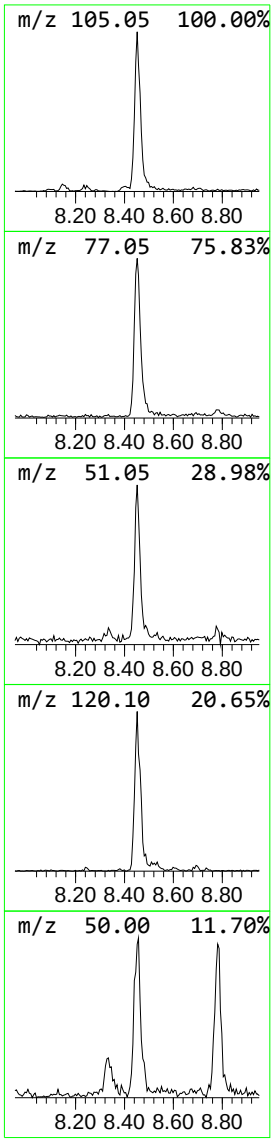
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Aceto phenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.452	3.43 ng/ul	100825	1,4-Dichlorobenzene-d4	7.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone	120	C8H8O	000098-86-2	93
2	Acetophenone	120	C8H8O	000098-86-2	91
3	Acetophenone	120	C8H8O	000098-86-2	90
4	Acetophenone	120	C8H8O	000098-86-2	90
5	Acetophenone	120	C8H8O	000098-86-2	86





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Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

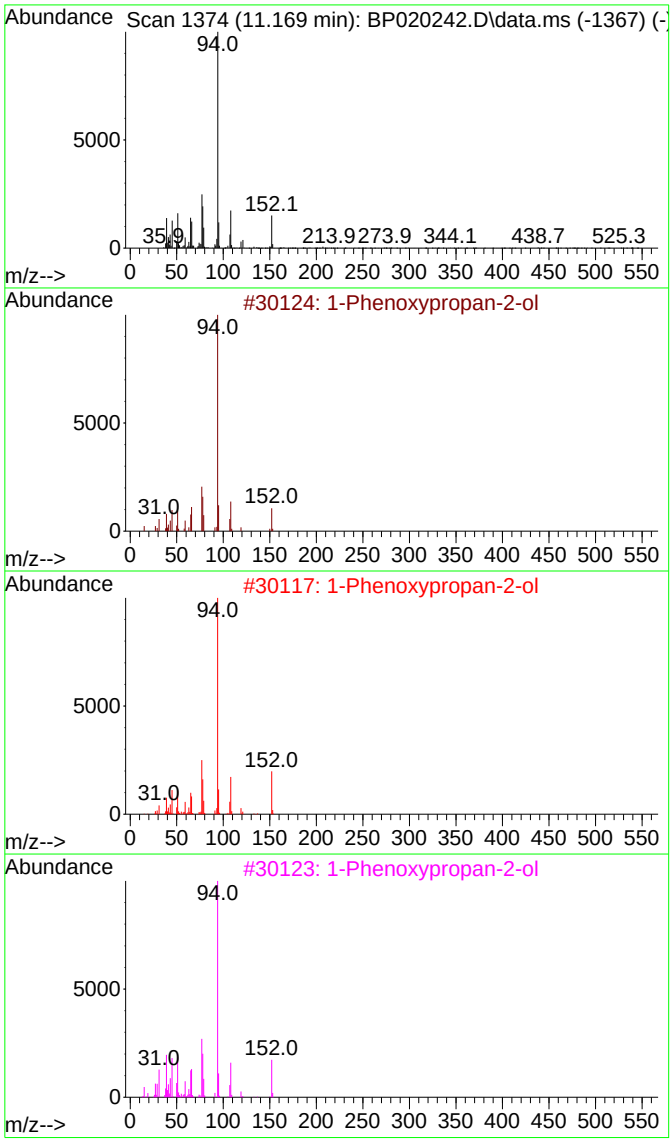
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

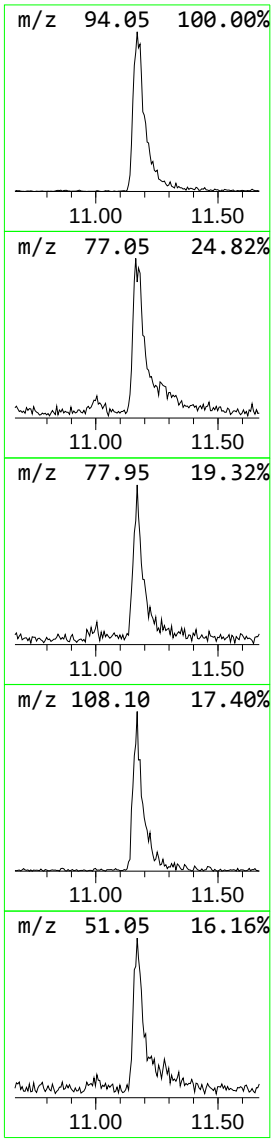
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.169	3.20 ng/ul	137778	Naphthalene-d8	10.387

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87





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Data File : BP020242.D
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

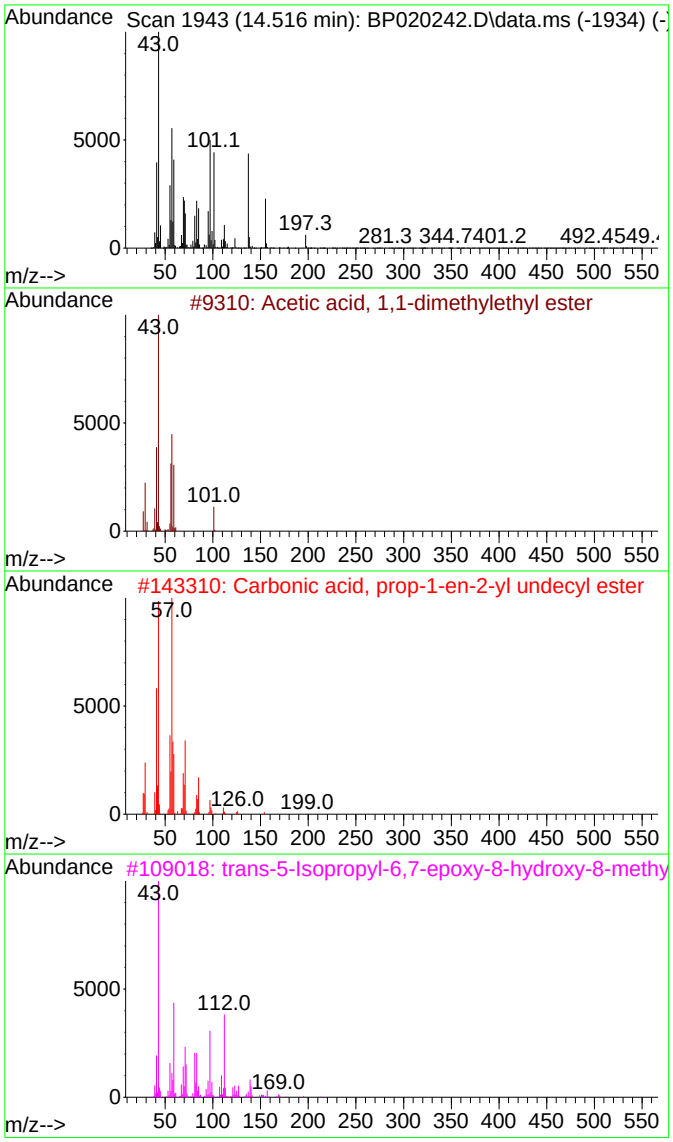
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

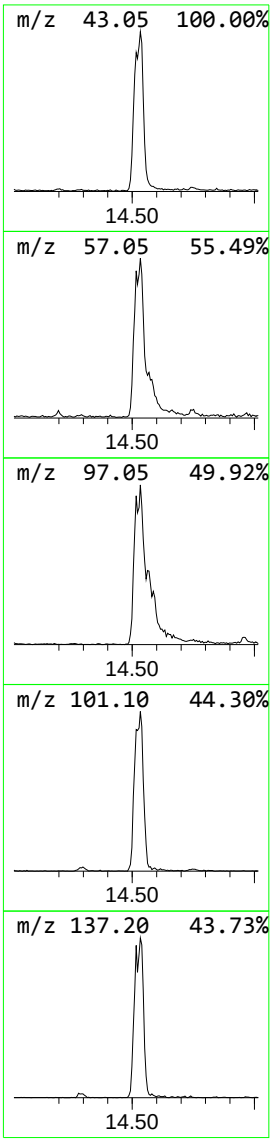
****

Peak Number 4 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.516	4.53 ng/ul	269170	Acenaphthene-d10	14.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	14
2	Carbonic acid, prop-1-en-2-yl un...	256	C15H28O3	1000382-90-6	10
3	trans-5-Isopropyl-6,7-epoxy-8-hy...	228	C13H24O3	058002-07-6	10
4	Pent-4-enoyl amide, 2-methyl-N-p...	155	C9H17NO	1000420-35-2	10
5	Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	10





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Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

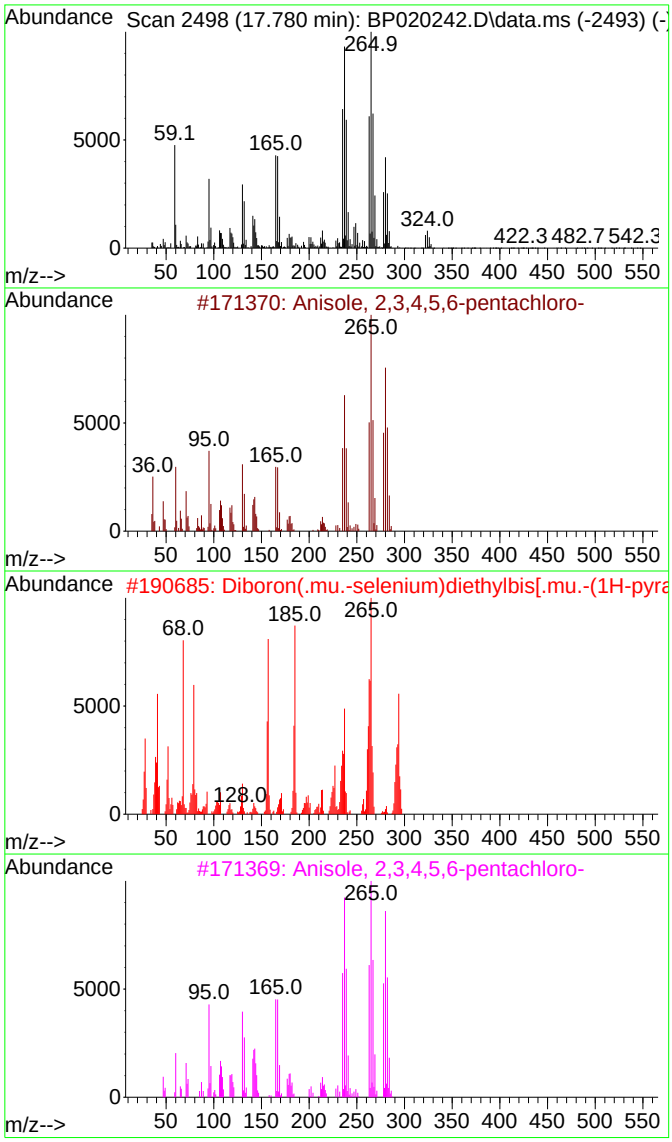
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

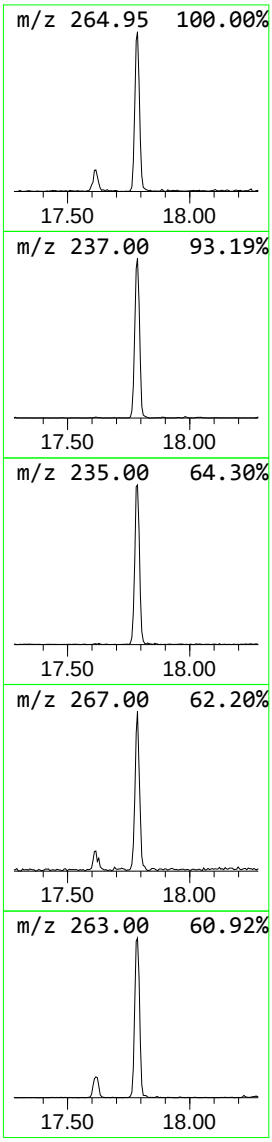
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Anisole, 2,3,4,5,6-pentachl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.781	3.39 ng/ul	250106	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	78
2	Diboron(.mu.-selenium)diethylbis...	294	C10H16B2N4Se	1000159-49-7	64
3	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	60
4	Rhodium, acetylanilinato-bis(eth...	293	C12H16NORh	1000153-76-6	30
5	Hexachlorocyclopentadiene	270	C5Cl6	000077-47-4	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

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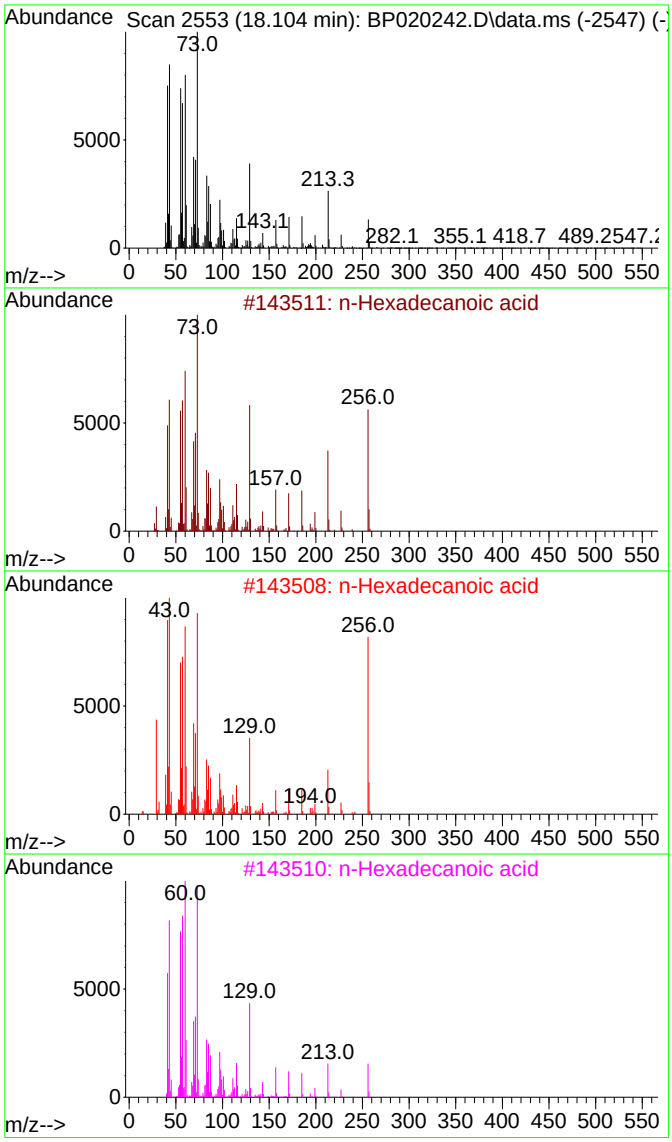
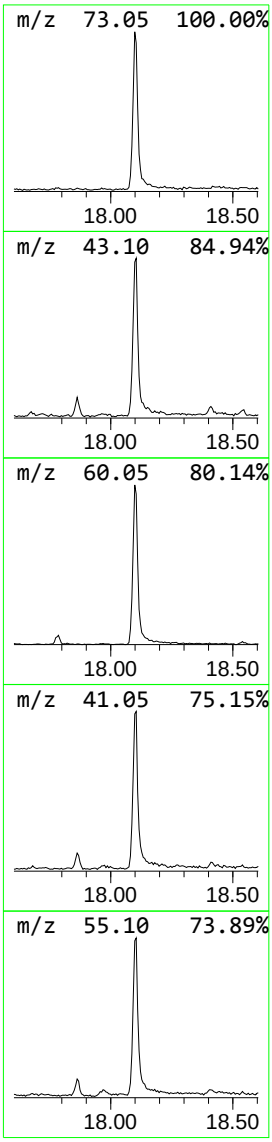
Peak Number 6 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	9.20 ng/ul	678890	Phenanthrene-d10	17.134

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	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

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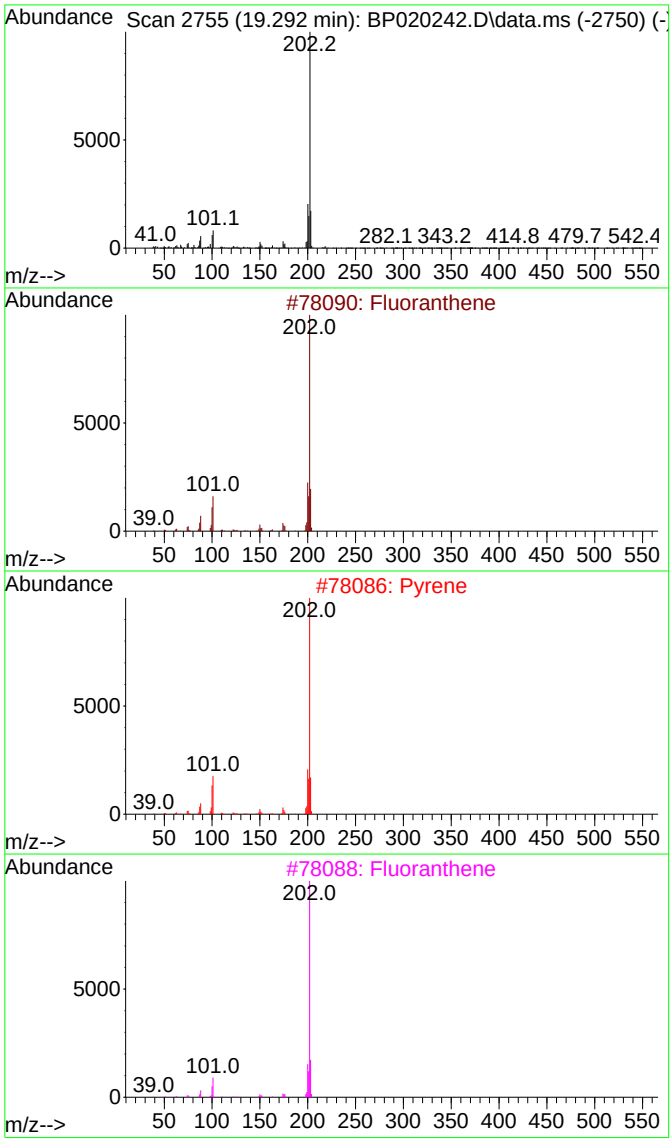
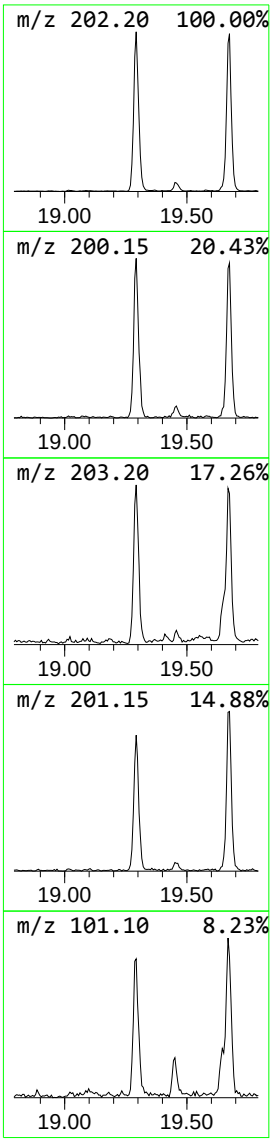
Peak Number 7 Fluoran thene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	3.81 ng/ul	281156	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene	202	C16H10	000206-44-0	91
2	Pyrene	202	C16H10	000129-00-0	90
3	Fluoranthene	202	C16H10	000206-44-0	81
4	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	81
5	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	81





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Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

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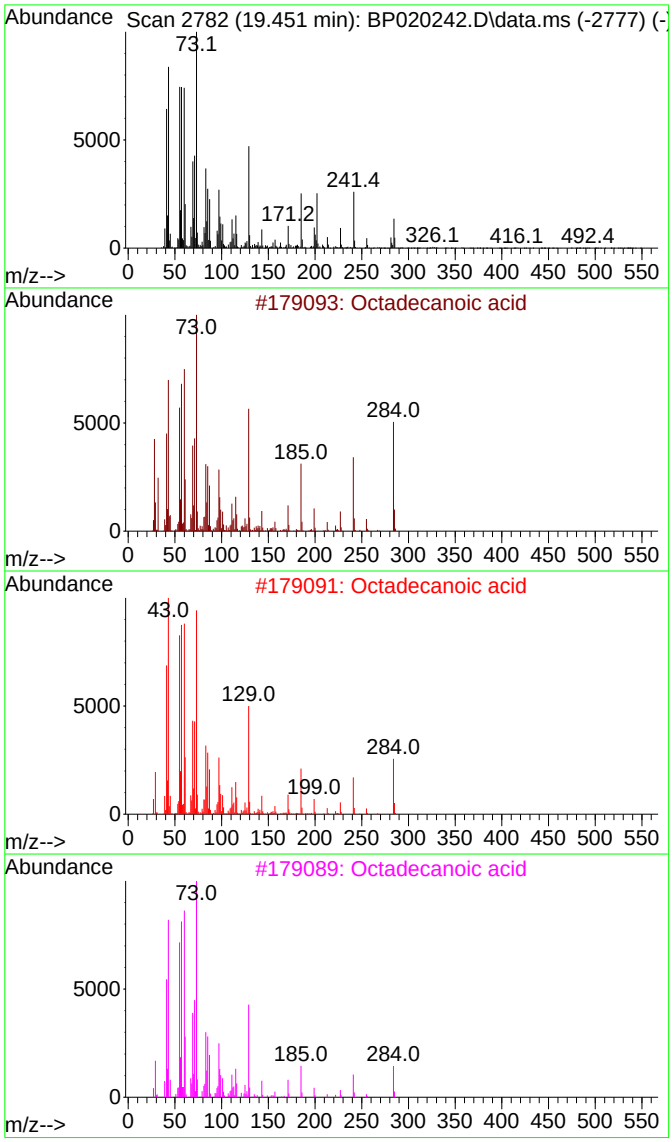
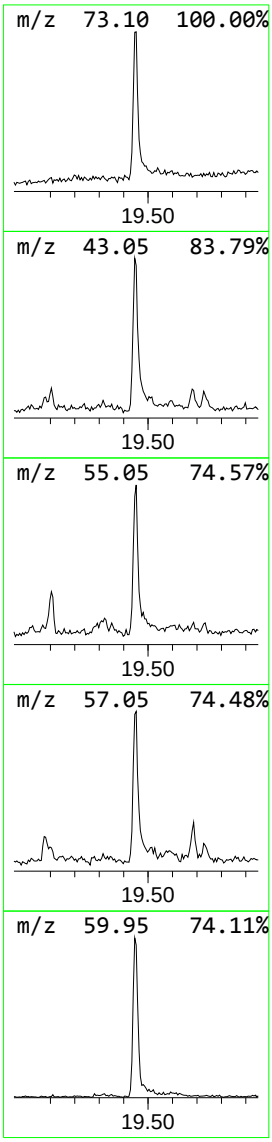
Peak Number 8 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	4.17 ng/ul	299342	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	96
4	Tridecanoic acid	214	C13H26O2	000638-53-9	70
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	68





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Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

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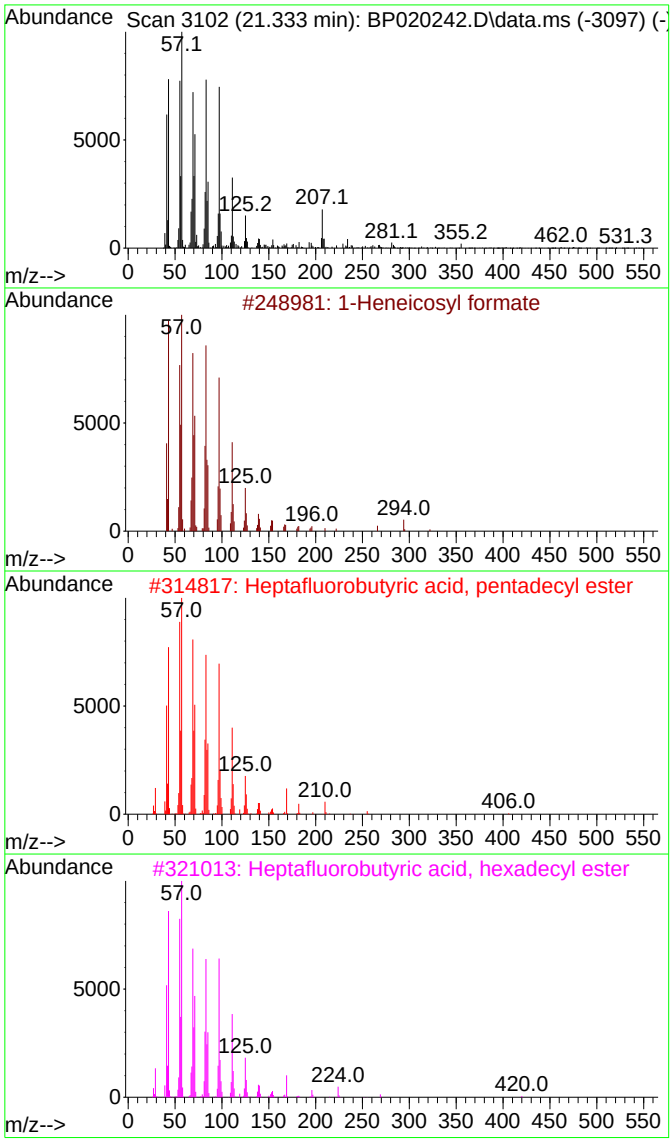
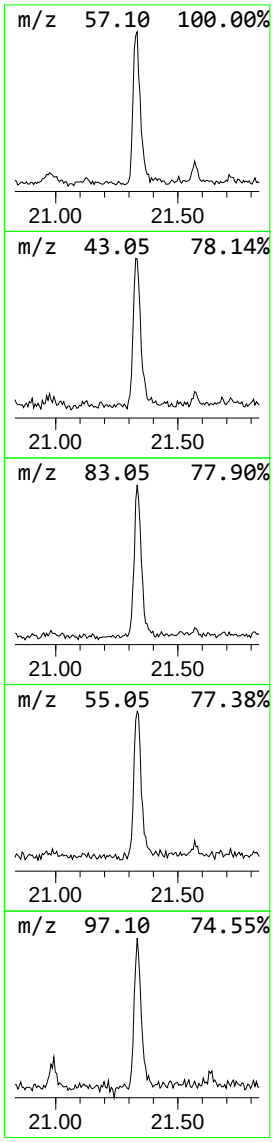
Peak Number 9 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.333	3.59 ng/ul	258100	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
2	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	90
4	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	90
5	1-Tricosene	322	C23H46	018835-32-0	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020242.D
Acq On : 08 May 2024 16:52
Operator : MA/JU
Sample : P2369-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.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
Propylene Glycol	3.481	3.4	ng/ul	99312	1	7.616	587452	20.0
Aceto phenone	8.452	3.4	ng/ul	100825	1	7.616	587452	20.0
1-Phenoxypropan...	11.169	3.2	ng/ul	137778	2	10.387	861437	20.0
unknown-01	14.516	4.5	ng/ul	269170	3	14.292	1189690	20.0
Anisole, 2,3,4,...	17.781	3.4	ng/ul	250106	4	17.134	1476430	20.0
n-Hexadecanoic ...	18.104	9.2	ng/ul	678890	4	17.134	1476430	20.0
Fluoran thene	19.292	3.8	ng/ul	281156	4	17.134	1476430	20.0
Octadecanoic acid	19.451	4.2	ng/ul	299342	5	21.633	1437300	20.0
1-Heneicosyl fo...	21.333	3.6	ng/ul	258100	5	21.633	1437300	20.0