LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

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-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012365.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	40	44	53	rVB	34269	56064	0.59%	0.065%
2	3.669	114	116	141	rBV	161629	310914	3.25%	0.362%
3	3.887	148	153	162	rVV	18433	30769	0.32%	0.036%
4	3.981	166	169	174	rVB	16415	20658	0.22%	0.024%
5	4.434	243	246	250	rVB3	8443	11691	0.12%	0.014%
6	4.593	270	273	279	rBV	18392	29563	0.31%	0.034%
7	4.916	322	328	341	rBV	1805951	2666520	27.89%	3.106%
8	5.558	433	437	445	rBV2	12565	22218	0.23%	0.026%
9	6.099	525	529	534	rBV4	8060	12824	0.13%	0.015%
10	6.852	652	657	663	rBV10	4708	11903	0.12%	0.014%
11	6.981	674	679	694	rBV2	83888	198229	2.07%	0.231%
12	7.140	699	706	720	rBV	564794	958040	10.02%	1.116%
13	7.334	732	739	760	rBV	308075	545065	5.70%	0.635%
14	7.805	812	819	827	rBV	832024	1347218	14.09%	1.569%
15	7.869	827	830	838	rVB2	19467	34025	0.36%	0.040%
16	8.522	930	941	958	rBV	299334	571124	5.97%	0.665%
17	8.910	1002	1007	1011	rBV5	7358	9872	0.10%	0.011%
18	8.975	1011	1018	1037	rBV	661488	1154420	12.08%	1.345%
19	9.128	1041	1044	1052	rVB9	5664	11784	0.12%	0.014%
20	9.357	1078	1083	1090	rVB	18817	33613	0.35%	0.039%
21	9.428	1090	1095	1106	rBV6	8646	18024	0.19%	0.021%
22	9.693	1134	1140	1151	rBV	277789	495221	5.18%	0.577%
23	9.999	1188	1192	1199	rVB4	9290	17864	0.19%	0.021%
24	10.234	1225	1232	1249	rBV	491192	909989	9.52%	1.060%
25	10.393	1252	1259	1285	rVV	558331	1269313	13.28%	1.478%
26	10.604	1288	1295	1310	rVV	1281119	2201226	23.03%	2.564%
27	10.751	1310	1320	1338	rVB	518616	1002227	10.48%	1.167%
28	11.281	1404	1410	1416	rBV4	7147	17945	0.19%	0.021%
29	11.822	1493	1502	1505	rBV7	6571	12814	0.13%	0.015%
30	11.963	1520	1526	1531	rVB2	12261	21215	0.22%	0.025%
31	12.022	1531	1536	1544	rVB10	7560	13030	0.14%	0.015%
32	12.140	1547	1556	1563	rBV	597553	1009620	10.56%	1.176%
33	12.198	1564	1566	1579	rVB3	25020	46230	0.48%	0.054%
34	12.804	1664	1669	1679	rBV	34742	74775	0.78%	0.087%
35	13.057	1706	1712	1720	rVB	61246	95323	1.00%	0.111%
36	13.269	1741	1748	1751	rBV	25667	40188	0.42%	0.047%

LSC Area Percent Report

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Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

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-BP102522.M
Title : SVOA CALIBRATION

37	13.345	1755	1761	1770	rVV	557152	929174	9.72%	1.082%
38	13.410	1770	1772	1778	rVB7	9411	16790	0.18%	0.020%
39	13.787	1830	1836	1843	rBV10	9488	27587	0.29%	0.032%
40	13.869	1843	1850	1864	rVV	2113626	2989933	31.28%	3.482%
41	14.145	1884	1897	1910	rBV	2242475	3490044	36.51%	4.065%
42	14.263	1915	1917	1922	rVV5	7866	11922	0.12%	0.014%
43	14.345	1926	1931	1936	rVV5	11990	19595	0.20%	0.023%
44	14.457	1942	1950	1959	rBV	2217169	3473149	36.33%	4.045%
45	14.528	1959	1962	1968	rVV4	19952	35766	0.37%	0.042%
46	14.692	1981	1990	2000	rBV4	45247	131487	1.38%	0.153%
47	14.857	2013	2018	2031	rBV4	35096	87097	0.91%	0.101%
48	14.957	2032	2035	2041	rVB2	12741	18357	0.19%	0.021%
49	15.210	2068	2078	2083	rBV	202502	301996	3.16%	0.352%
50	15.281	2087	2090	2100	rVB2	48827	85134	0.89%	0.099%
51	15.386	2104	2108	2111	rBV2	13246	20630	0.22%	0.024%
52	15.457	2111	2120	2131	rBV2	6085038	9559271	100.00%	11.134%
53	15.586	2136	2142	2153	rBV	470548	698360	7.31%	0.813%
54	15.681	2154	2158	2161	rBV2	36228	52703	0.55%	0.061%
55	15.804	2175	2179	2186	rBV	38469	57551	0.60%	0.067%
56	15.910	2192	2197	2203	rBV3	48759	103399	1.08%	0.120%
57	15.975	2203	2208	2212	rVB	104627	145259	1.52%	0.169%
58	16.198	2243	2246	2249	rVV	21168	26517	0.28%	0.031%
59	16.292	2257	2262	2266	rBV	71864	107739	1.13%	0.125%
60	16.351	2267	2272	2278	rVB3	101125	186253	1.95%	0.217%
61	16.416	2278	2283	2287	rBV2	72378	107987	1.13%	0.126%
62	16.457	2287	2290	2295	rVB2	37995	48887	0.51%	0.057%
63	16.510	2295	2299	2305	rBV2	31027	71217	0.75%	0.083%
64	16.616	2312	2317	2324	rVB4	68212	126015	1.32%	0.147%
65	16.681	2324	2328	2332	rBV	65740	93595	0.98%	0.109%
66	16.751	2337	2340	2346	rVB2	42309	65811	0.69%	0.077%
67	16.810	2348	2350	2355	rVB4	7365	10296	0.11%	0.012%
68	16.928	2367	2370	2375	rVB	31644	43097	0.45%	0.050%
69	16.969	2375	2377	2380	rBV3	10115	10751	0.11%	0.013%
70	17.216	2412	2419	2430	rBV	3113913	4403657	46.07%	5.129%
71	17.316	2430	2436	2445	rBV	4541006	6347586	66.40%	7.393%
72	17.616	2483	2487	2497	rVB9	27664	58765	0.61%	0.068%
73	17.757	2507	2511	2514	rVB5	15946	22301	0.23%	0.026%
74	17.886	2528	2533	2541	rBV	293168	388223	4.06%	0.452%
75	18.104	2565	2570	2579	rBV	553691	884682	9.25%	1.030%
76	18.386	2614	2618	2629	rBV3	57134	102777	1.08%	0.120%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

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-BP102522.M
Title : SVOA CALIBRATION

77	18.510	2637	2639	2643	rVB4	23048	21604	0.23%	0.025%
78	18.945	2708	2713	2718	rBV2	116773	222565	2.33%	0.259%
79	18.986	2718	2720	2721	rVV2	68599	60657	0.63%	0.071%
80	19.016	2721	2725	2736	rVB	1144939	1434878	15.01%	1.671%
81	19.145	2741	2747	2752	rBV	479290	583666	6.11%	0.680%
82	19.198	2752	2756	2759	rBV	137382	199293	2.08%	0.232%
83	19.333	2776	2779	2793	rBV	234945	376048	3.93%	0.438%
84	19.557	2811	2817	2822	rBV	6290847	8151593	85.27%	9.494%
85	19.792	2853	2857	2862	rBV2	177635	274532	2.87%	0.320%
86	20.057	2898	2902	2908	rBV3	390863	719248	7.52%	0.838%
87	20.169	2918	2921	2925	rVB	158991	179077	1.87%	0.209%
88	20.245	2930	2934	2938	rVV	145015	181975	1.90%	0.212%
89	20.398	2956	2960	2965	rBV	887318	979190	10.24%	1.140%
90	20.463	2968	2971	2974	rBV	125432	155070	1.62%	0.181%
91	20.516	2978	2980	2983	rVV	143332	142901	1.49%	0.166%
92	20.551	2983	2986	2992	rVB2	159088	209399	2.19%	0.244%
93	21.021	3061	3066	3074	rBV2	575151	1034867	10.83%	1.205%
94	21.310	3110	3115	3125	rBV	4528418	5782612	60.49%	6.735%
95	21.392	3127	3129	3133	rVB	164880	189279	1.98%	0.220%
96	23.551	3489	3496	3509	rVB2	4291607	8832326	92.40%	10.287%
97	23.704	3515	3522	3537	rVB	2719328	5486766	57.40%	6.390%

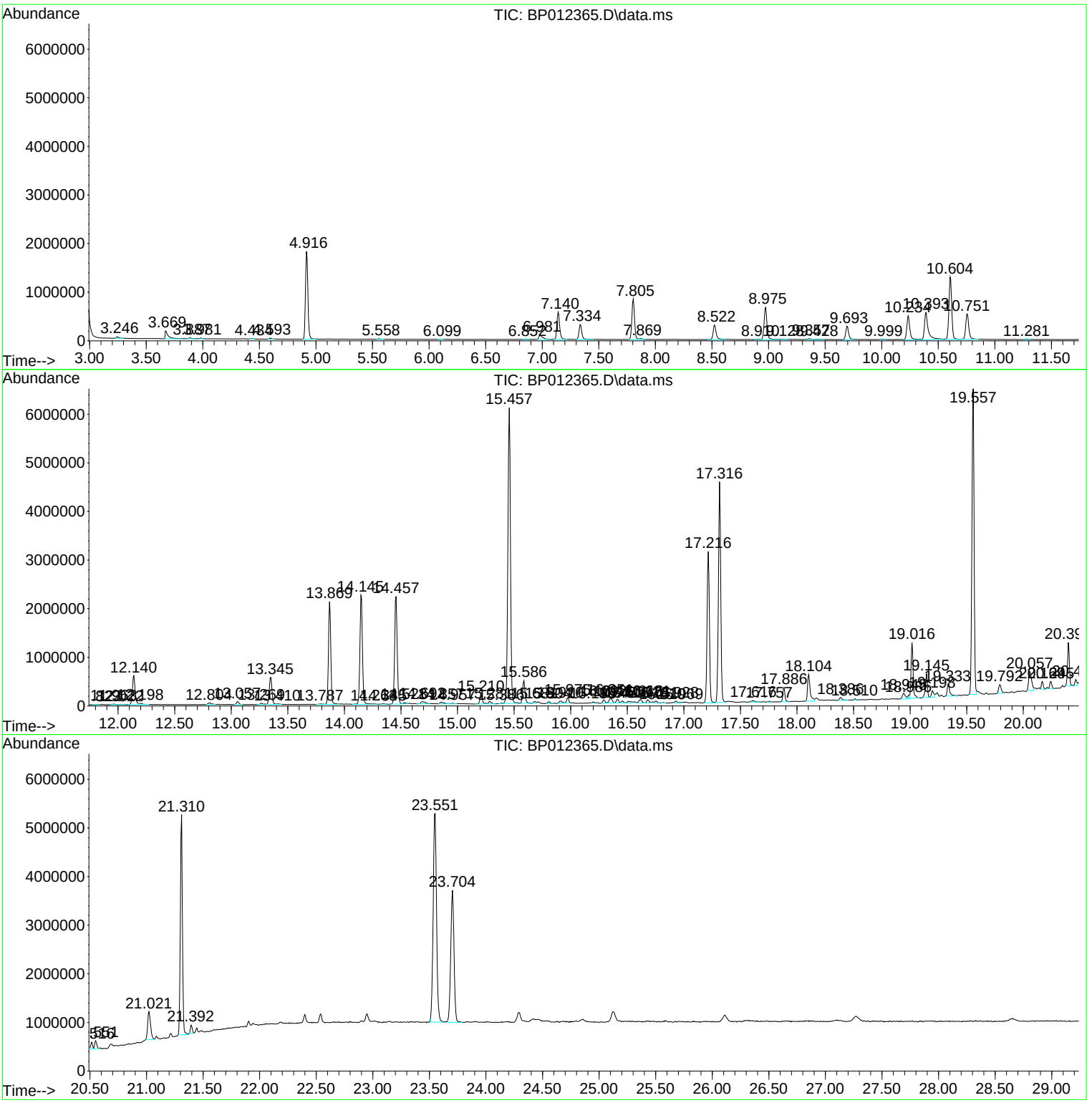
Sum of corrected areas: 85860424

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

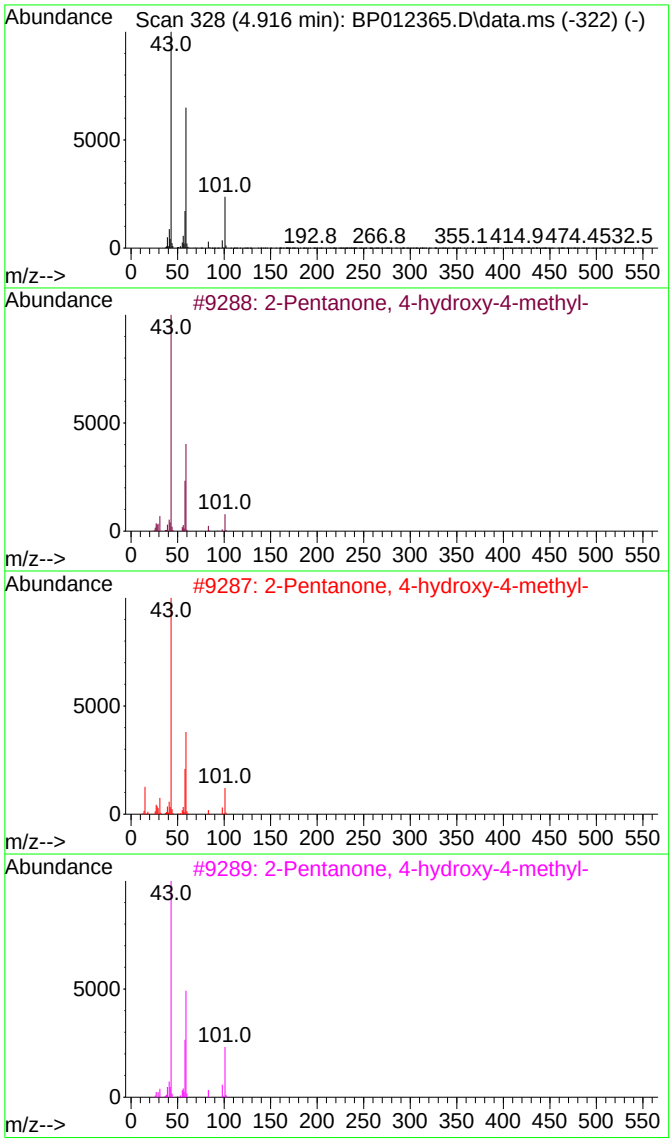
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

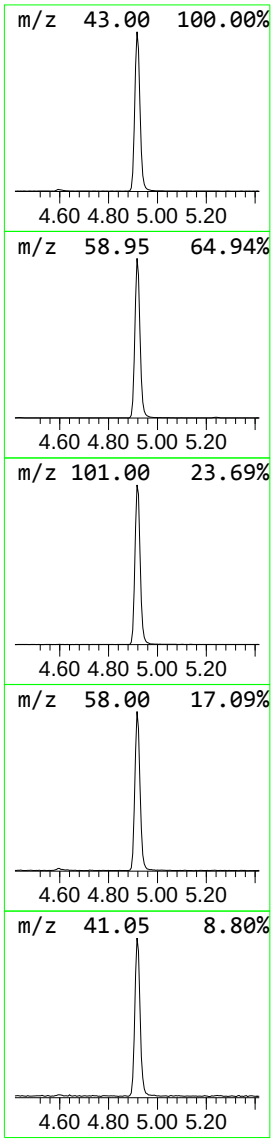
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.916	39.59 ng/ul	2666520	1,4-Dichlorobenzene-d4	7.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

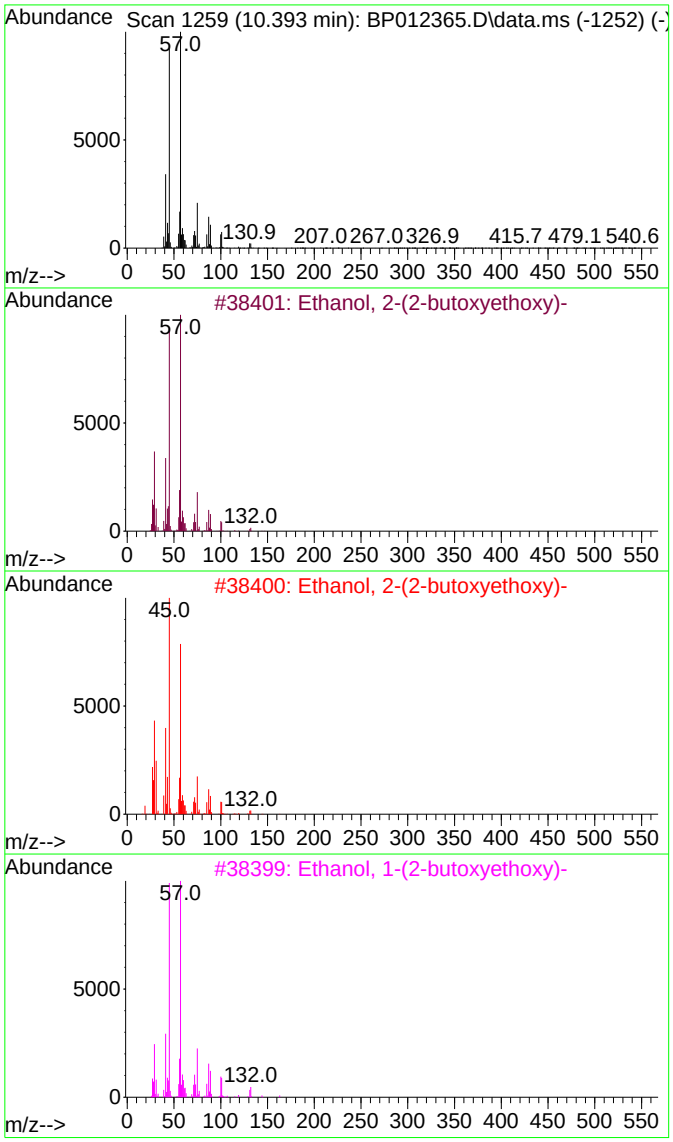
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

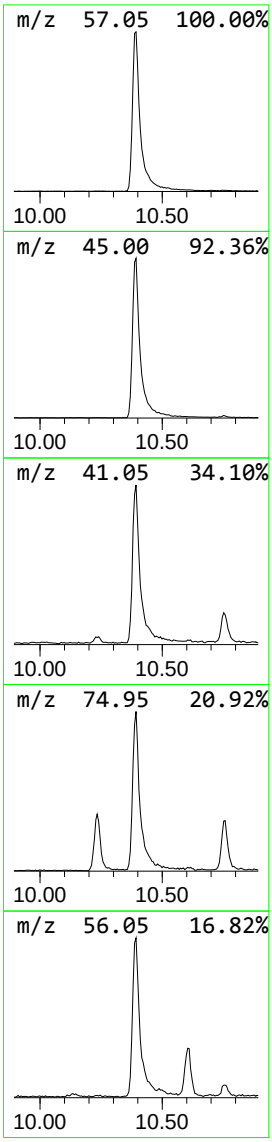
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.393	11.53 ng/ul	1269310	Naphthalene-d8	10.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

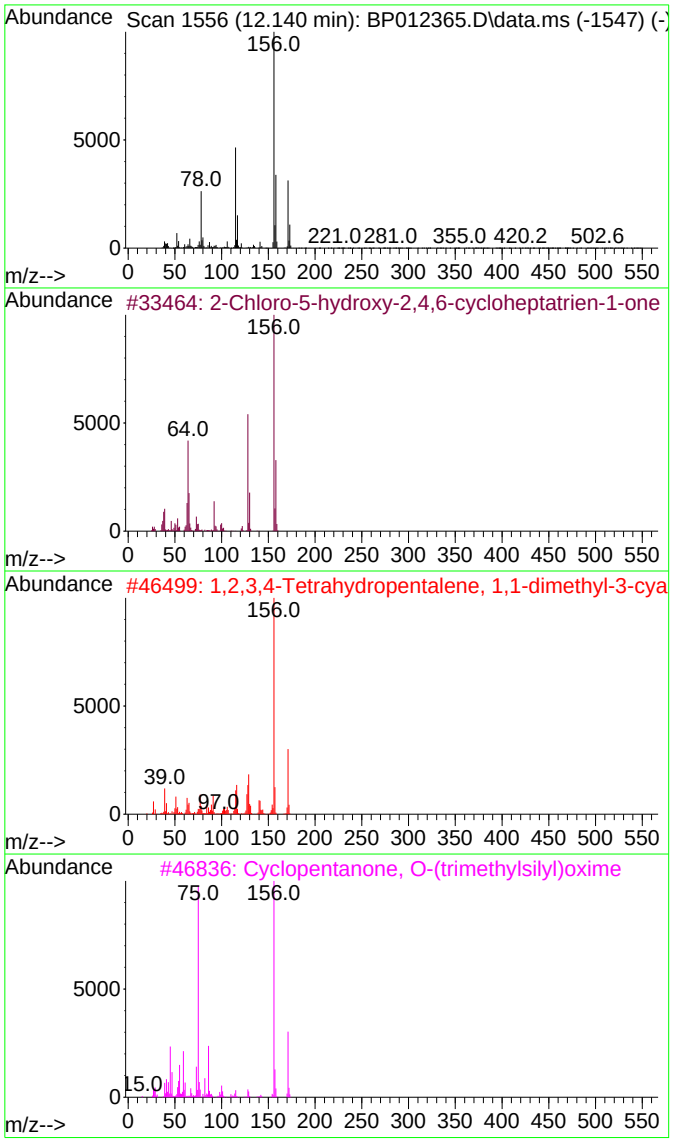
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

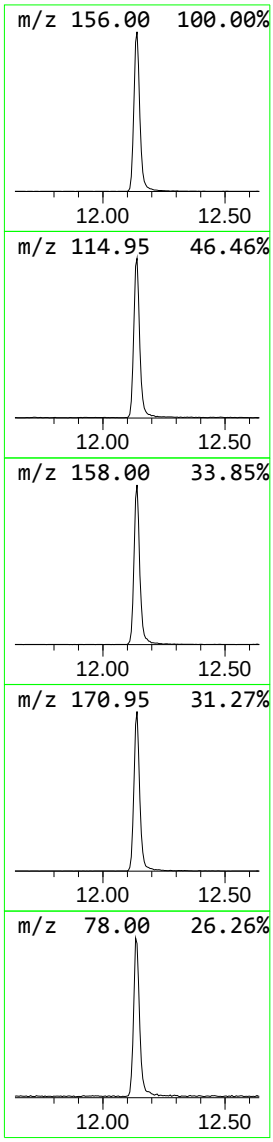
****

Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.140	9.17 ng/ul	1009620	Naphthalene-d8	10.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37
3	Cyclopentanone, O-(trimethylsily...	171	C8H17NOSi	026208-87-7	32
4	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
5	Chloroxylenol	156	C8H9ClO	000088-04-0	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
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Instrument :
BNA_P
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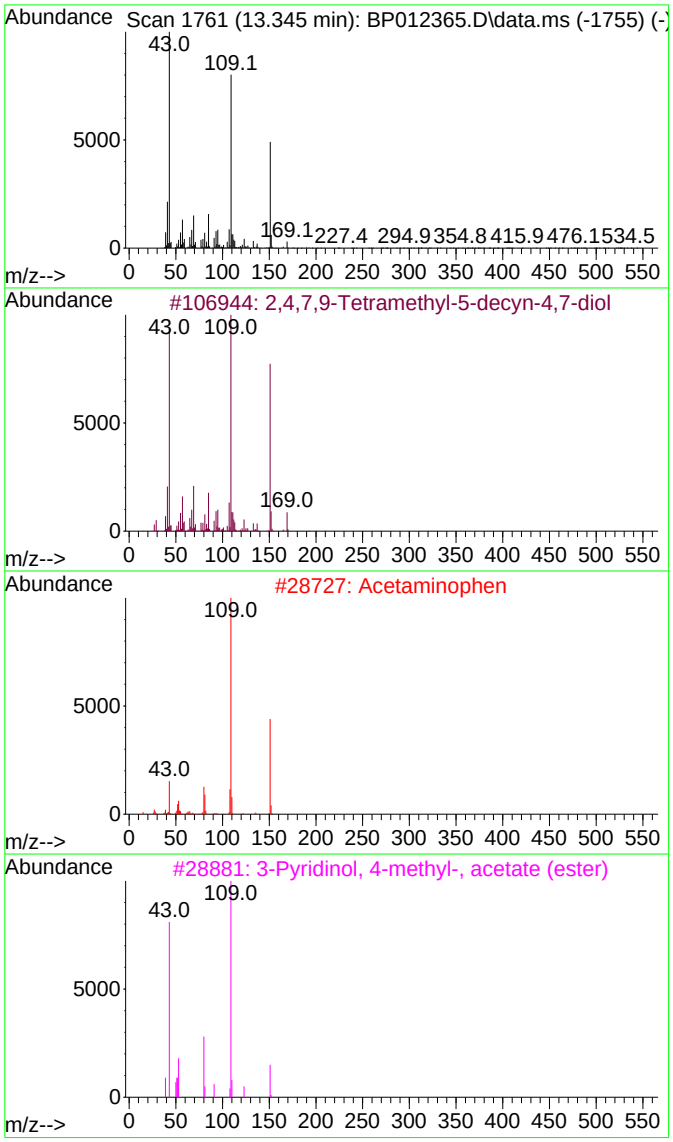
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

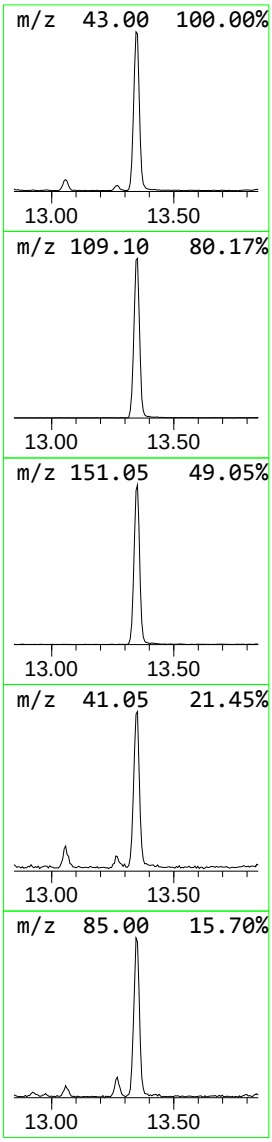
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.345	5.35 ng/ul	929174	Acenaphthene-d10	14.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	Acetaminophen	151	C8H9NO2	000103-90-2	59
3	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59
4	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
5	Metacetamol	151	C8H9NO2	000621-42-1	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

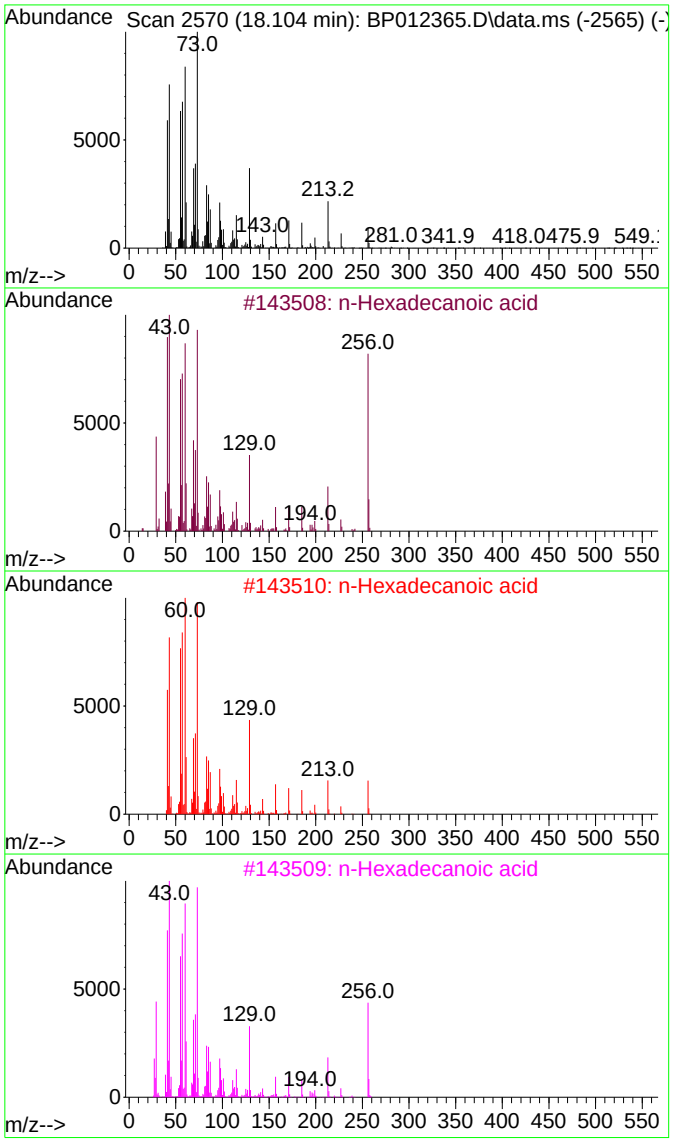
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

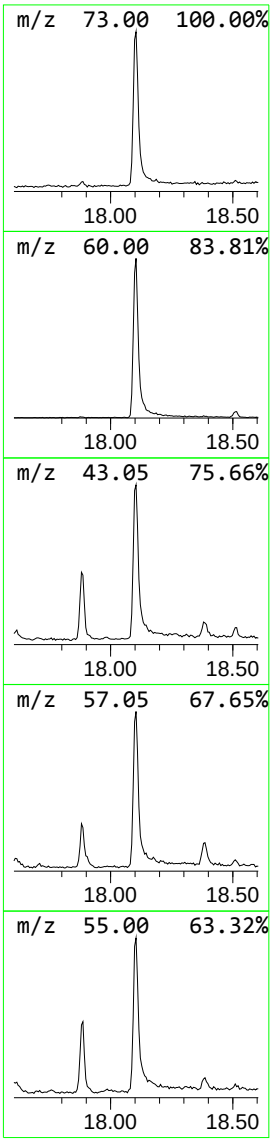
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	4.02 ng/ul	884682	Phenanthrene-d10	17.216

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	97
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95





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Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

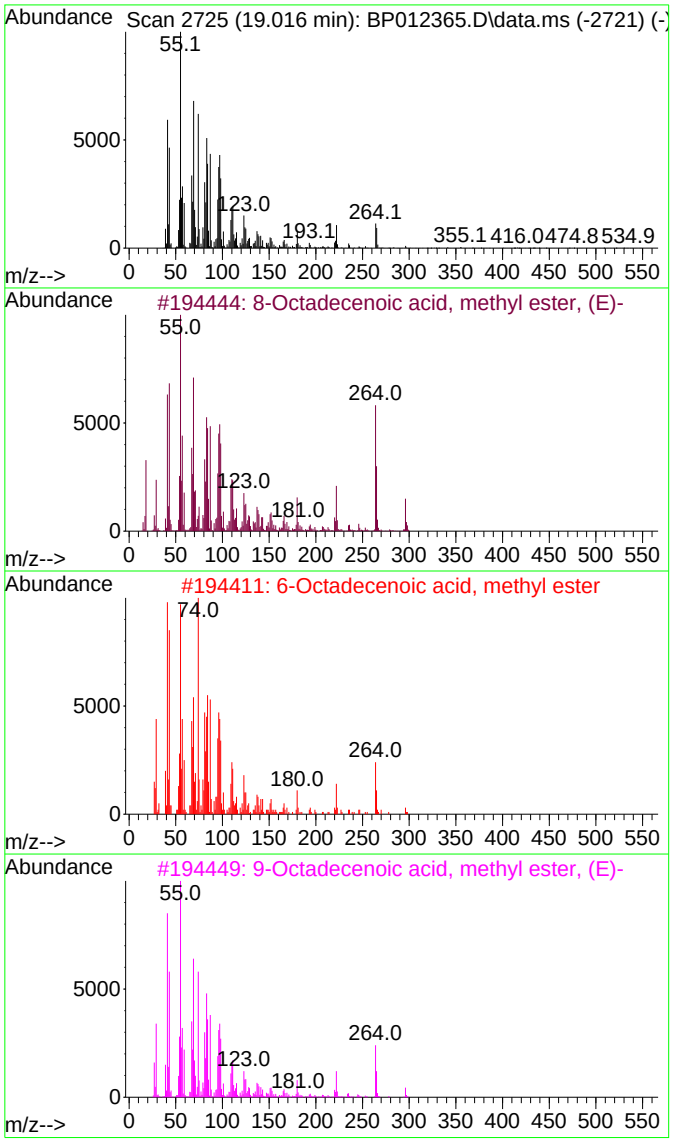
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

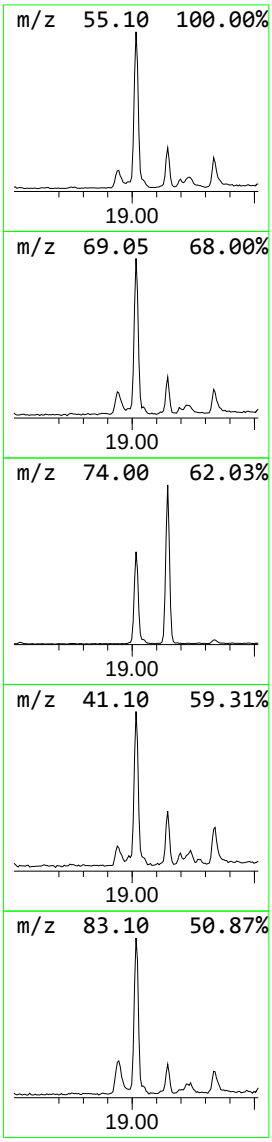
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 8-Octadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.016	6.52 ng/ul	1434880	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	99
2	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
3	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
4	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
5	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

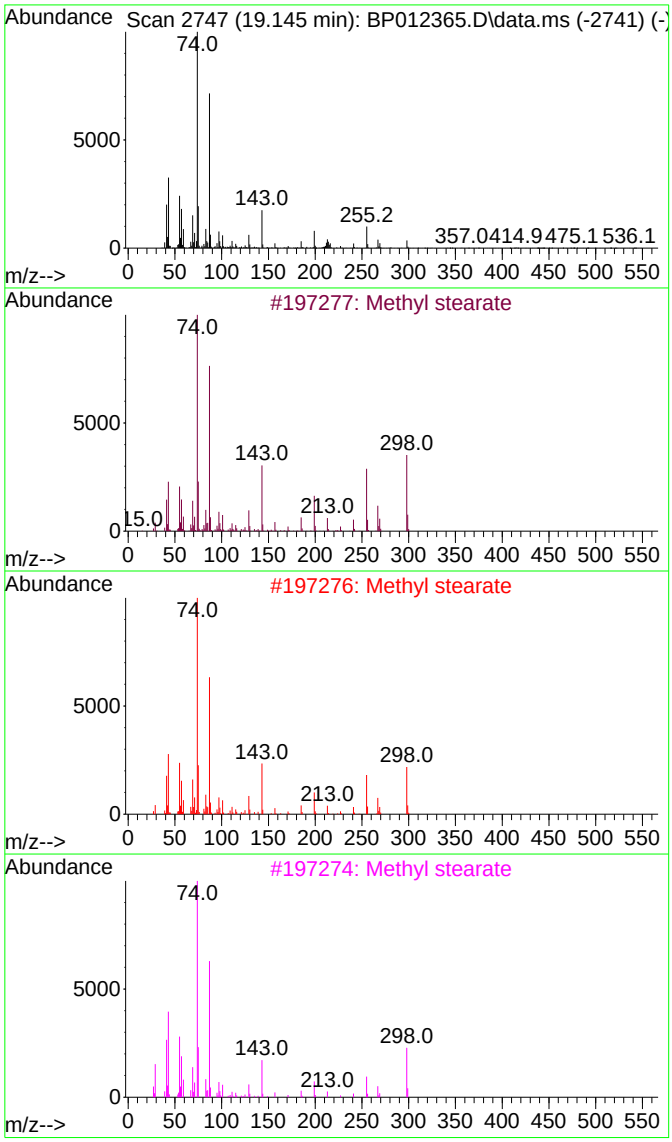
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

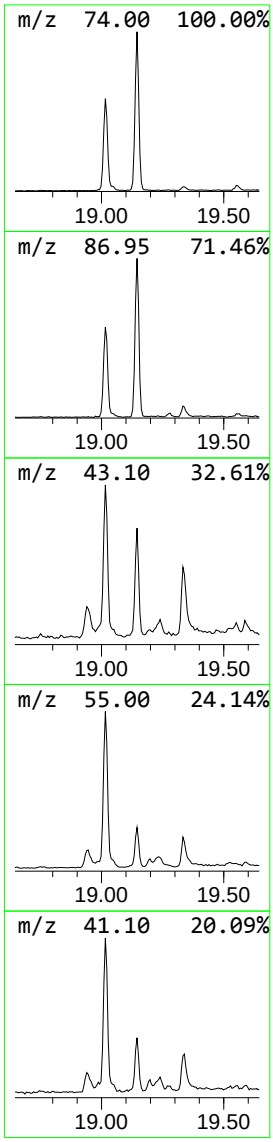
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Methyl stearate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.145	2.65 ng/ul	583666	Phenanthrene-d10	17.216

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	97
2	Methyl stearate	298	C19H38O2	000112-61-8	97
3	Methyl stearate	298	C19H38O2	000112-61-8	97
4	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
5	Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	91





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Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

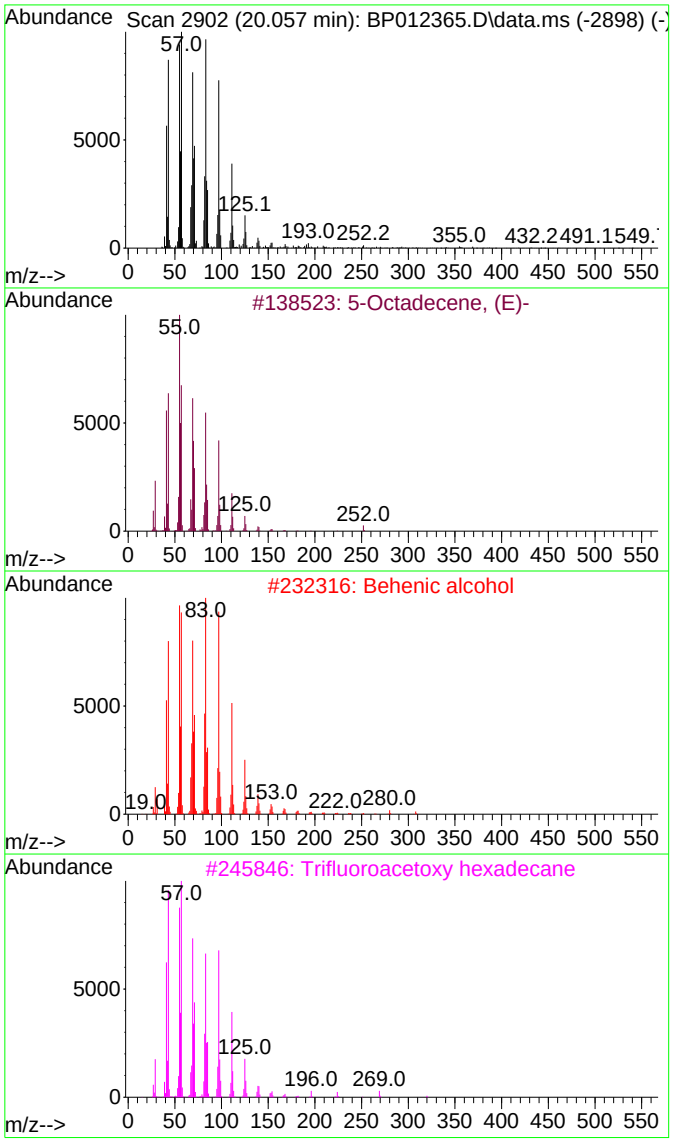
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

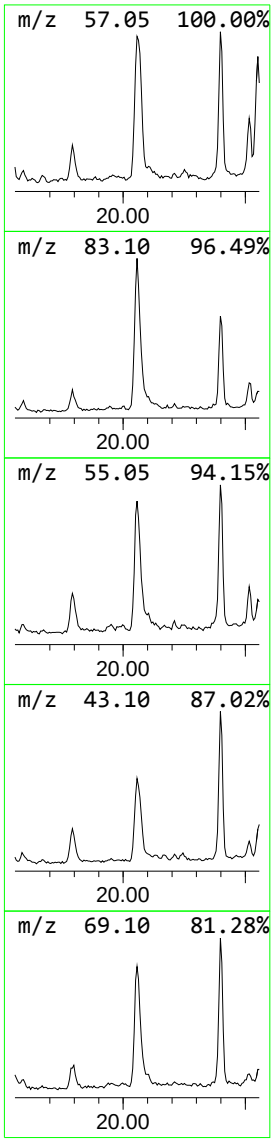
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.057	2.49 ng/ul	719248	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5	1-Octadecanol	270	C18H38O	000112-92-5	91





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Data File : BP012365.D
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Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

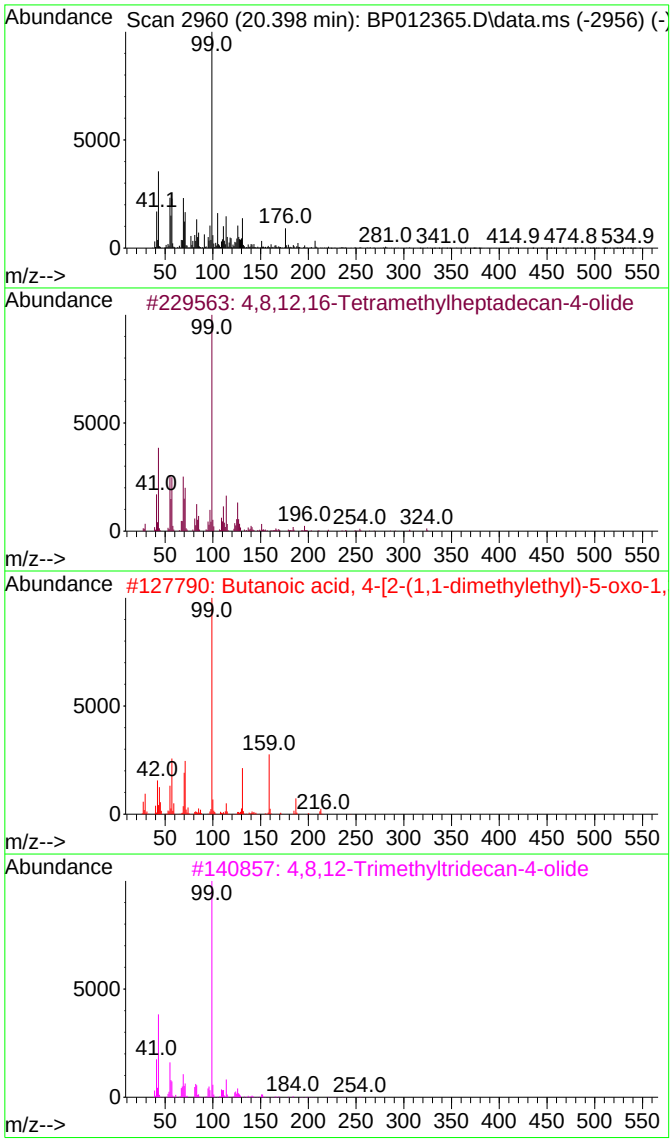
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

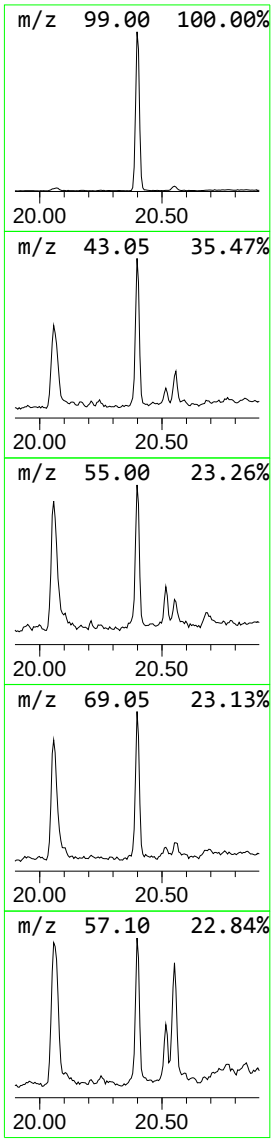
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 4,8,12,16-Tetramethylheptad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.398	3.39 ng/ul	979190	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	72
2	Butanoic acid, 4-[2-(1,1-dimethy...	244	C12H20O5	1000196-81-8	50
3	4,8,12-Trimethyltridecan-4-olide	254	C16H30O2	220904-24-5	49
4	4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	49
5	2H-Pyran-2-one, tetrahydro-6-pen...	170	C10H18O2	000705-86-2	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

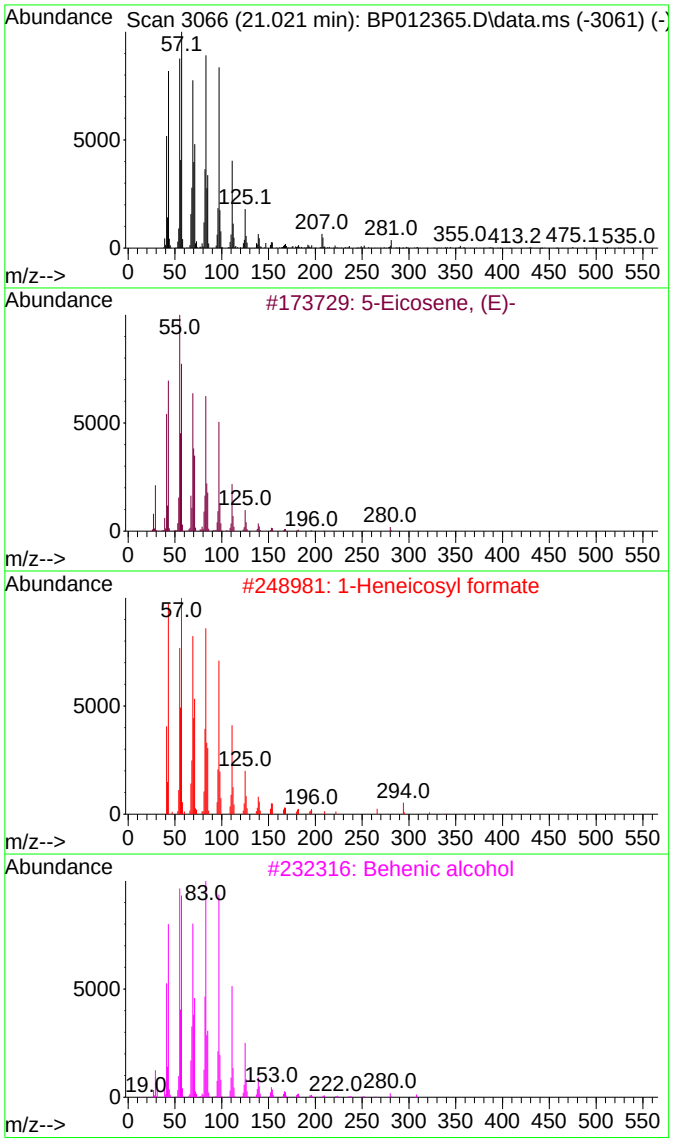
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

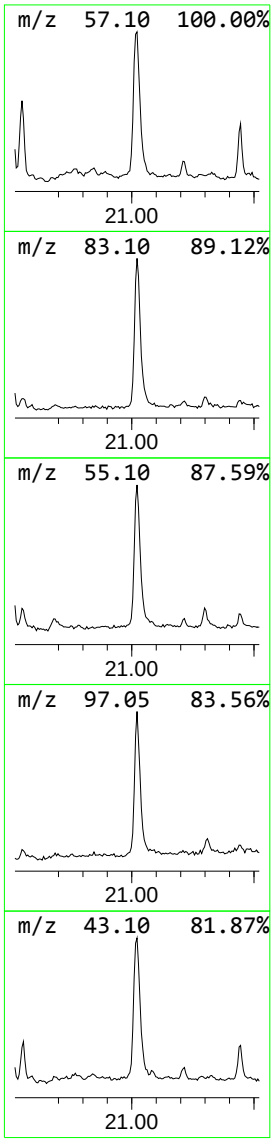
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	3.58 ng/ul	1034870	Chrysene-d12	21.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
3	Behenic alcohol	326	C22H46O	000661-19-8	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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
2-Pentanone, 4-...	4.916	39.6	ng/ul	2666520	1	7.805	1347220	20.0
Ethanol, 2-(2-b...	10.393	11.5	ng/ul	1269310	2	10.604	2201230	20.0
unknown-01	12.140	9.2	ng/ul	1009620	2	10.604	2201230	20.0
2,4,7,9-Tetrame...	13.345	5.3	ng/ul	929174	3	14.457	3473150	20.0
n-Hexadecanoic ...	18.104	4.0	ng/ul	884682	4	17.216	4403660	20.0
8-Octadecenoic ...	19.016	6.5	ng/ul	1434880	4	17.216	4403660	20.0
Methyl stearate	19.145	2.6	ng/ul	583666	4	17.216	4403660	20.0
(DEL) Alkane: S...	20.057	2.5	ng/ul	719248	5	21.310	5782610	20.0
4,8,12,16-Tetra...	20.398	3.4	ng/ul	979190	5	21.310	5782610	20.0
(DEL) Alkane: S...	21.022	3.6	ng/ul	1034870	5	21.310	5782610	20.0