LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

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-BP101922.M
Title : SVOA CALIBRATION

Signal : TIC: BP012265.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	7	9	29	rVB	364443	512696	8.25%	0.473%
2	3.258	42	46	54	rVB	87325	113601	1.83%	0.105%
3	3.681	115	118	131	rBV	428076	629136	10.12%	0.581%
4	3.781	131	135	142	rVB	433989	546852	8.80%	0.505%
5	3.899	151	155	162	rVB2	102420	153651	2.47%	0.142%
6	3.999	167	172	180	rVB	122439	180131	2.90%	0.166%
7	4.558	262	267	274	rVB	97753	143821	2.31%	0.133%
8	4.917	324	328	330	rBV	59633	91278	1.47%	0.084%
9	5.022	339	346	358	rVB2	70508	154034	2.48%	0.142%
10	5.452	413	419	427	rBV	115973	208479	3.35%	0.192%
11	5.822	475	482	488	rBV	190304	284800	4.58%	0.263%
12	6.299	555	563	570	rBV3	64386	130733	2.10%	0.121%
13	6.793	639	647	652	rBV	56818	101355	1.63%	0.094%
14	6.993	670	681	695	rVB	1247458	2296871	36.96%	2.121%
15	7.158	702	709	724	rVB	1167847	1976285	31.80%	1.825%
16	7.352	735	742	751	rBV	1381368	2238152	36.01%	2.066%
17	7.428	751	755	757	rVV	61627	84097	1.35%	0.078%
18	7.458	757	760	767	rVB	109658	171029	2.75%	0.158%
19	7.816	810	821	831	rVV	1601573	2720384	43.77%	2.512%
20	7.928	835	840	847	rVV2	57635	114201	1.84%	0.105%
21	8.534	936	943	953	rBV	839139	1432989	23.06%	1.323%
22	8.987	1013	1020	1026	rVV	1335887	2231522	35.91%	2.060%
23	9.040	1026	1029	1038	rVB	74558	108248	1.74%	0.100%
24	9.152	1043	1048	1057	rVB8	31654	83053	1.34%	0.077%
25	9.710	1133	1143	1154	rBV	1118238	1920362	30.90%	1.773%
26	10.246	1228	1234	1243	rVV	1637396	2798231	45.03%	2.583%
27	10.404	1255	1261	1276	rBV	527850	1016425	16.36%	0.938%
28	10.622	1287	1298	1303	rBV	2274883	3961828	63.75%	3.658%
29	10.675	1303	1307	1314	rVB	322267	543202	8.74%	0.502%
30	10.769	1316	1323	1341	rVB	730240	1422739	22.89%	1.314%
31	11.640	1466	1471	1483	rVV	79133	154909	2.49%	0.143%
32	12.040	1533	1539	1549	rBV	101652	161205	2.59%	0.149%
33	12.287	1575	1581	1589	rVB	435795	719220	11.57%	0.664%
34	12.504	1613	1618	1627	rVV	367864	592961	9.54%	0.547%
35	12.993	1695	1701	1708	rVB2	64432	104569	1.68%	0.097%
36	13.287	1744	1751	1758	rBV2	171867	332208	5.35%	0.307%

LSC Area Percent Report

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Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

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-BP101922.M
Title : SVOA CALIBRATION

37	13.493	1782	1786	1789	rBV	51301	87465	1.41%	0.081%
38	13.628	1804	1809	1810	rBV	91746	118282	1.90%	0.109%
39	13.787	1830	1836	1841	rVV	236598	409908	6.60%	0.378%
40	13.840	1841	1845	1848	rVV	199761	307525	4.95%	0.284%
41	13.887	1848	1853	1859	rVV	2755401	3990182	64.20%	3.684%
42	13.940	1859	1862	1867	rVV	120718	165995	2.67%	0.153%
43	14.022	1872	1876	1880	rVV	137593	202521	3.26%	0.187%
44	14.057	1880	1882	1887	rVB2	71763	87606	1.41%	0.081%
45	14.163	1893	1900	1917	rVB	3357794	5394706	86.80%	4.981%
46	14.357	1928	1933	1941	rVB	161613	217968	3.51%	0.201%
47	14.469	1942	1952	1959	rBV	3321768	4957065	79.76%	4.577%
48	14.687	1982	1989	1997	rBV	714612	1309898	21.08%	1.209%
49	14.869	2011	2020	2029	rBV3	193122	471609	7.59%	0.435%
50	14.945	2029	2033	2036	rVB	62862	85598	1.38%	0.079%
51	14.987	2036	2040	2048	rVB6	42449	105401	1.70%	0.097%
52	15.087	2053	2057	2062	rBV	70500	112644	1.81%	0.104%
53	15.139	2062	2066	2071	rVB	63378	84080	1.35%	0.078%
54	15.257	2079	2086	2090	rBV4	111658	226188	3.64%	0.209%
55	15.316	2090	2096	2103	rVB2	187363	299069	4.81%	0.276%
56	15.469	2115	2122	2127	rVV	4155340	6112211	98.35%	5.643%
57	15.510	2127	2129	2135	rVV4	207961	311124	5.01%	0.287%
58	15.592	2138	2143	2154	rVB	690338	1012611	16.29%	0.935%
59	15.722	2160	2165	2174	rVB4	101434	211690	3.41%	0.195%
60	15.816	2174	2181	2187	rVV	122537	214122	3.45%	0.198%
61	15.963	2202	2206	2213	rVV	128790	255664	4.11%	0.236%
62	16.057	2215	2222	2229	rVB3	51925	121660	1.96%	0.112%
63	16.204	2238	2247	2253	rBV	429642	922324	14.84%	0.852%
64	16.763	2335	2342	2346	rBV3	138698	288056	4.64%	0.266%
65	16.881	2358	2362	2369	rBV2	94498	153960	2.48%	0.142%
66	16.981	2371	2379	2383	rVV2	273892	444932	7.16%	0.411%
67	17.034	2384	2388	2398	rVB4	96264	233027	3.75%	0.215%
68	17.122	2399	2403	2411	rBV	148813	255366	4.11%	0.236%
69	17.192	2411	2415	2416	rBV2	80718	98223	1.58%	0.091%
70	17.228	2416	2421	2426	rVV	3269655	4840872	77.89%	4.469%
71	17.269	2426	2428	2433	rVV	577420	753326	12.12%	0.695%
72	17.328	2433	2438	2449	rVB2	3809335	5755441	92.61%	5.314%
73	17.722	2501	2505	2509	rBV	268402	305212	4.91%	0.282%
74	17.898	2531	2535	2541	rVB	168444	217854	3.51%	0.201%
75	18.004	2548	2553	2559	rBV3	181003	317866	5.11%	0.293%
76	18.063	2559	2563	2566	rBV	123494	175450	2.82%	0.162%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

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-BP101922.M
Title : SVOA CALIBRATION

77	18.116	2566	2572	2576	rBV	812726	1321829	21.27%	1.220%
78	18.281	2596	2600	2604	rBV3	156366	262771	4.23%	0.243%
79	18.322	2604	2607	2612	rVB	140955	189475	3.05%	0.175%
80	18.398	2615	2620	2625	rBV2	390996	610781	9.83%	0.564%
81	18.586	2648	2652	2656	rBV	108255	147805	2.38%	0.136%
82	19.004	2716	2723	2726	rBV2	293263	527418	8.49%	0.487%
83	19.028	2726	2727	2731	rVB2	129389	127187	2.05%	0.117%
84	19.128	2740	2744	2753	rVB4	103370	232322	3.74%	0.214%
85	19.210	2755	2758	2759	rBV	129114	127441	2.05%	0.118%
86	19.239	2759	2763	2772	rVB	1290140	1939403	31.21%	1.791%
87	19.351	2778	2782	2786	rBV2	333122	462417	7.44%	0.427%
88	19.569	2814	2819	2822	rBV	4530813	6214761	100.00%	5.738%
89	19.663	2833	2835	2841	rVB2	112235	158060	2.54%	0.146%
90	20.080	2904	2906	2911	rVB2	313874	362886	5.84%	0.335%
91	20.569	2986	2989	2993	rBV	217122	237694	3.82%	0.219%
92	20.810	3028	3030	3034	rVB	183891	201029	3.23%	0.186%
93	21.027	3063	3067	3076	rVB4	834538	1500347	24.14%	1.385%
94	21.322	3111	3117	3121	rBV2	3729399	5606480	90.21%	5.176%
95	21.357	3121	3123	3129	rVB	653765	778790	12.53%	0.719%
96	22.992	3394	3401	3403	rBV2	934200	2099034	33.77%	1.938%
97	23.027	3405	3407	3419	rVB2	1381214	2480827	39.92%	2.290%
98	23.568	3493	3499	3504	rBV2	2714130	5259336	84.63%	4.856%
99	23.721	3520	3525	3531	rVB2	2152956	4084819	65.73%	3.771%
100	24.315	3623	3626	3634	rVB	675048	1311571	21.10%	1.211%

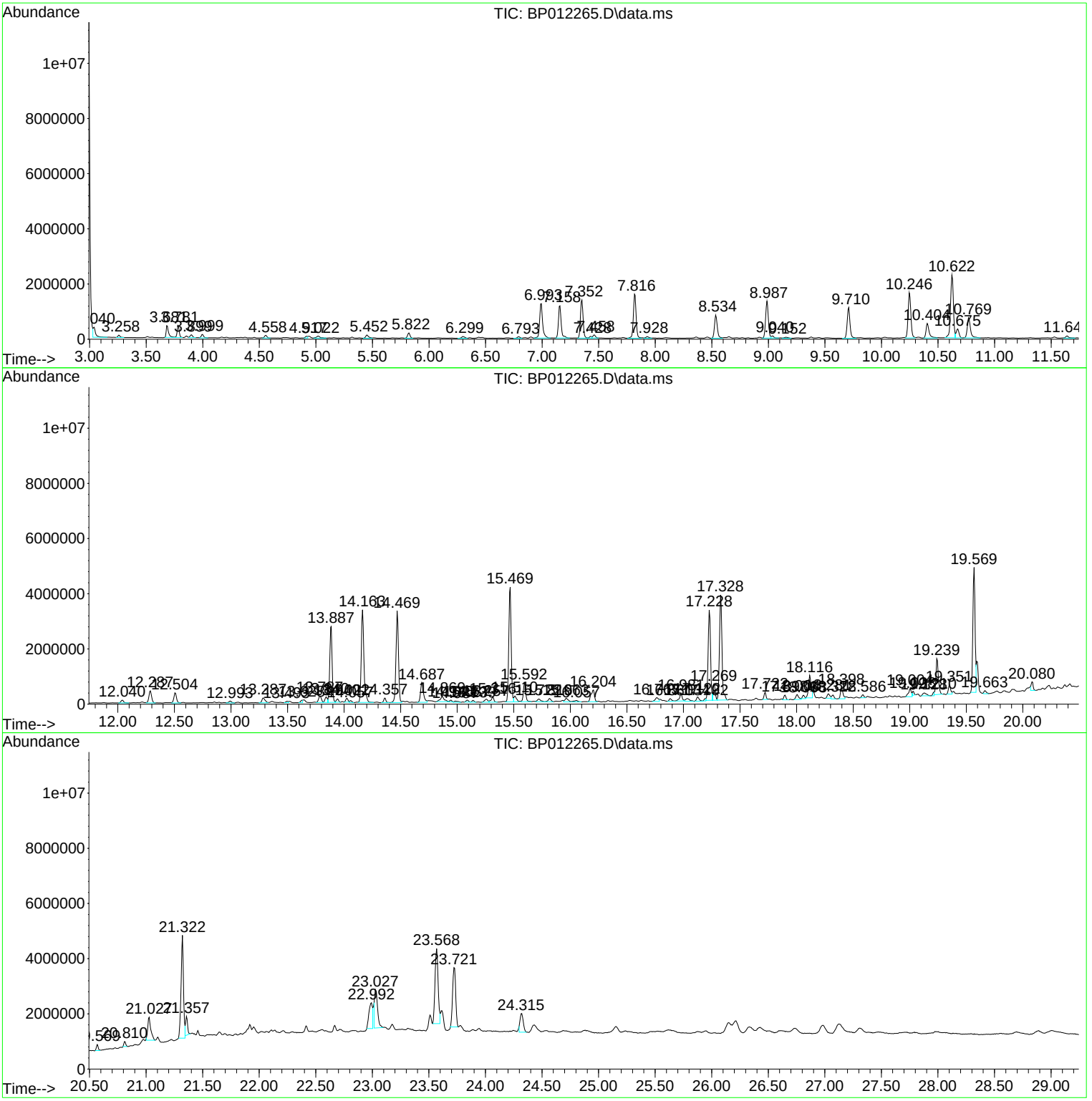
Sum of corrected areas: 108314441

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
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Sample : N5064-03
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ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

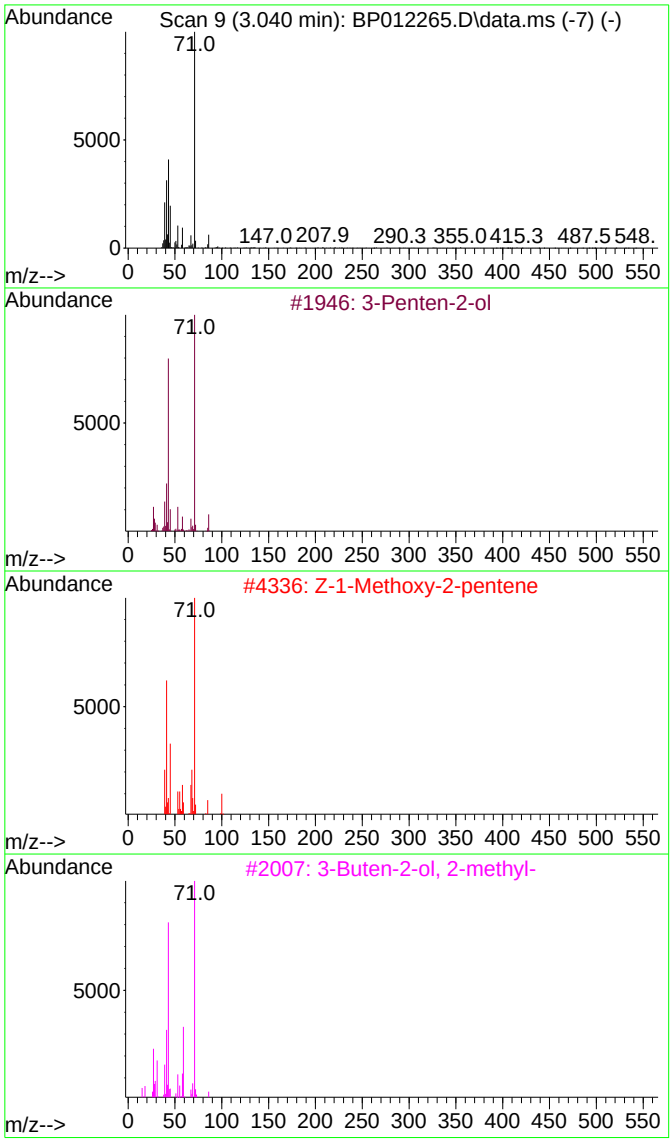
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

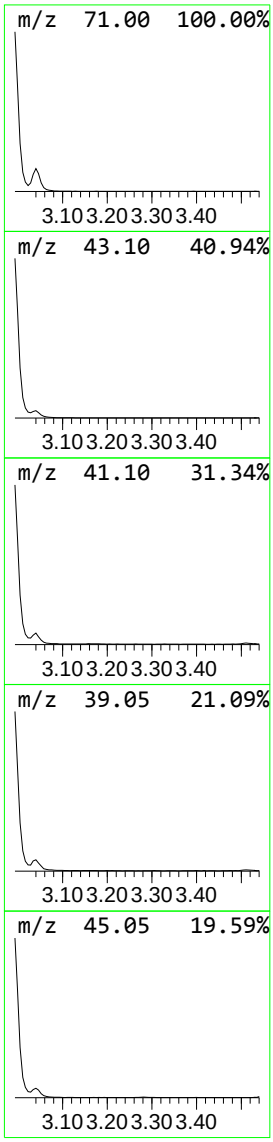
****

Peak Number 1 3-Penten-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	3.77 ng/ul	512696	1,4-Dichlorobenzene-d4	7.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-ol	86	C5H10O	001569-50-2	72
2	Z-1-Methoxy-2-pentene	100	C6H12O	1000279-59-4	64
3	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	59
4	Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	59
5	2-Propenoic acid, 2-methyl-, (te...	170	C9H14O3	002455-24-5	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

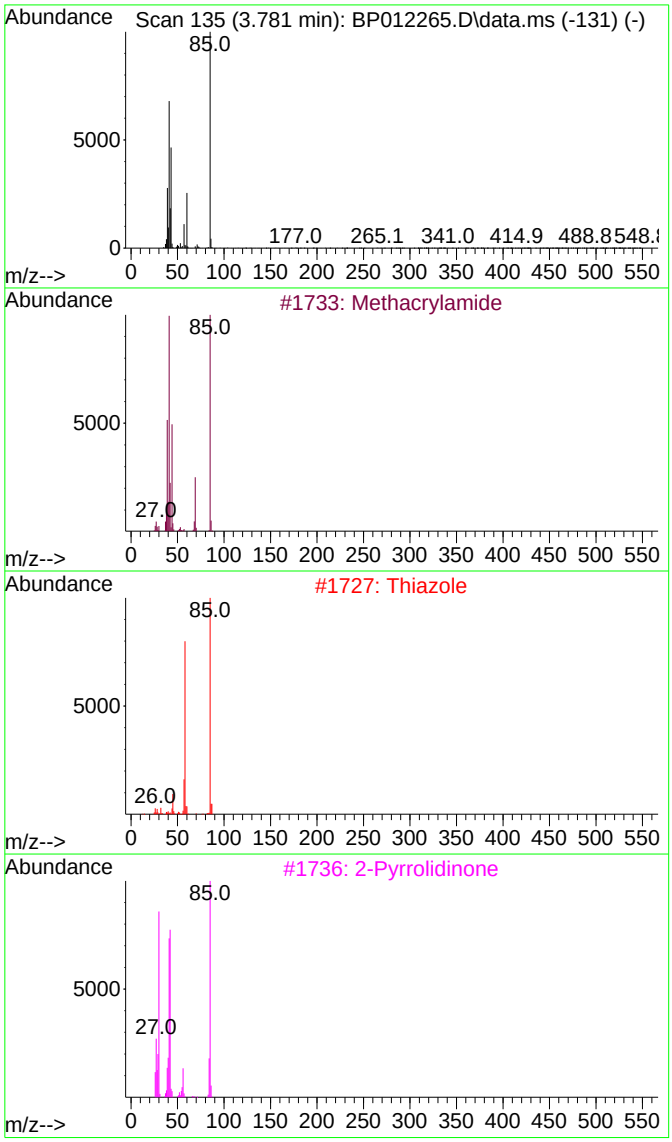
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

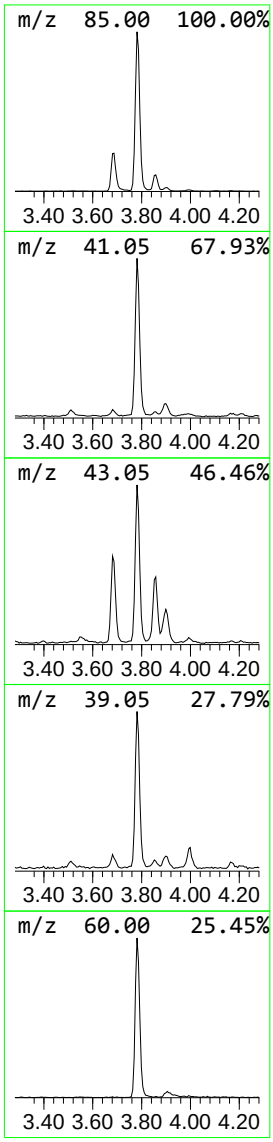
****

Peak Number 2 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.781	4.02 ng/ul	546852	1,4-Dichlorobenzene-d4	7.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methacrylamide	85	C4H7NO	000079-39-0	33
2	Thiazole	85	C3H3NS	000288-47-1	28
3	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
4	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

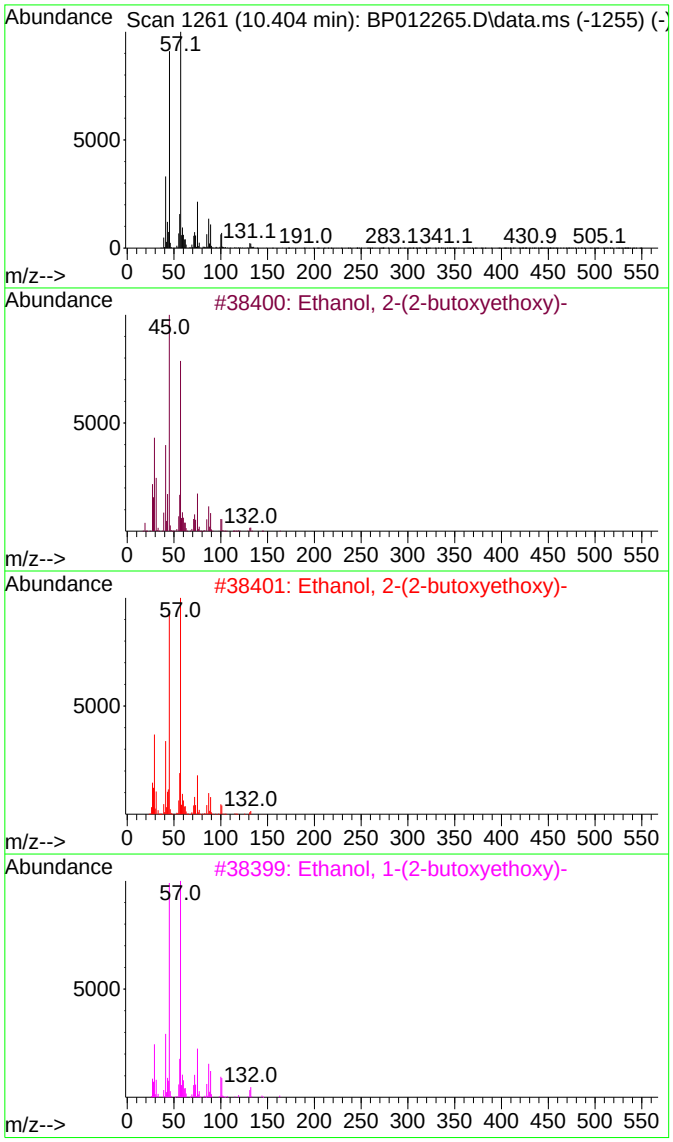
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

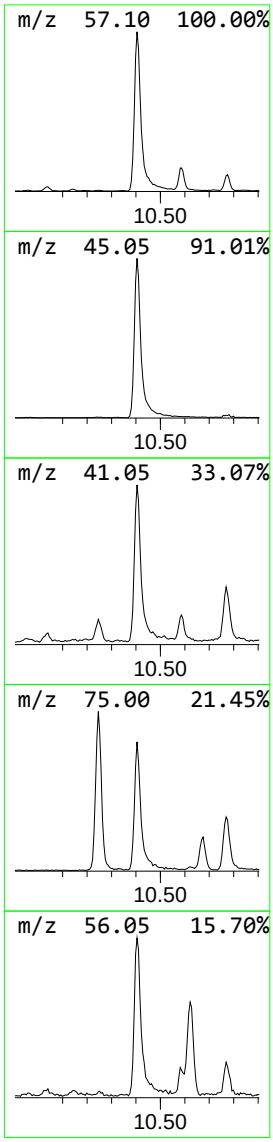
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.404	5.13 ng/ul	1016430	Naphthalene-d8	10.622

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	83
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

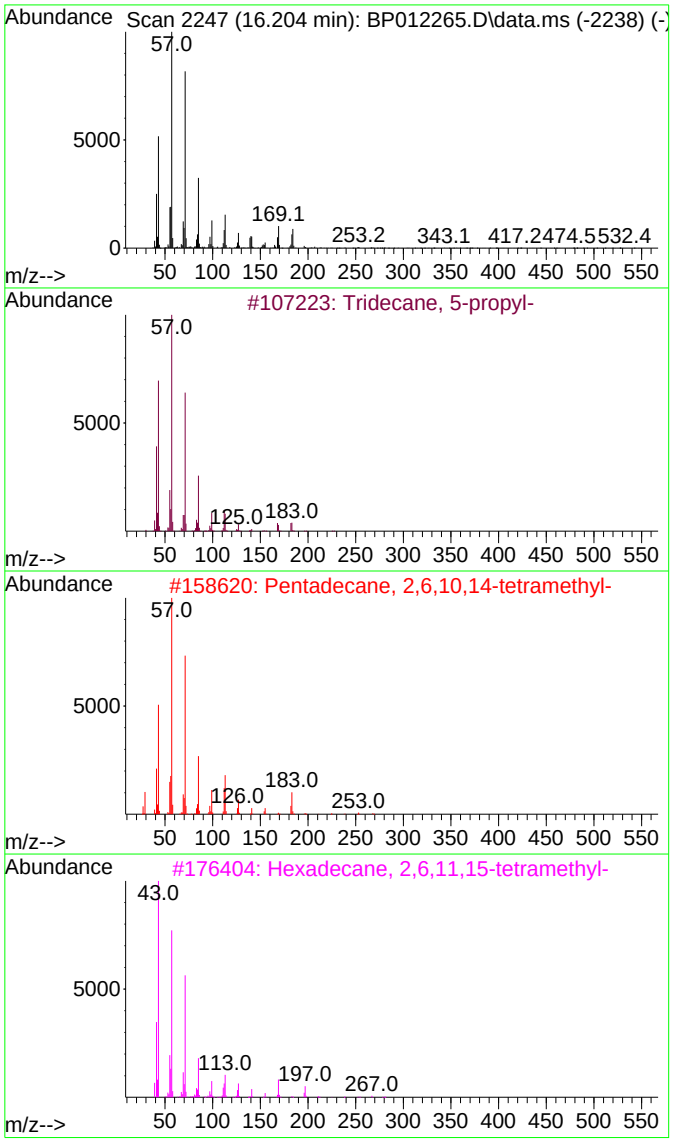
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

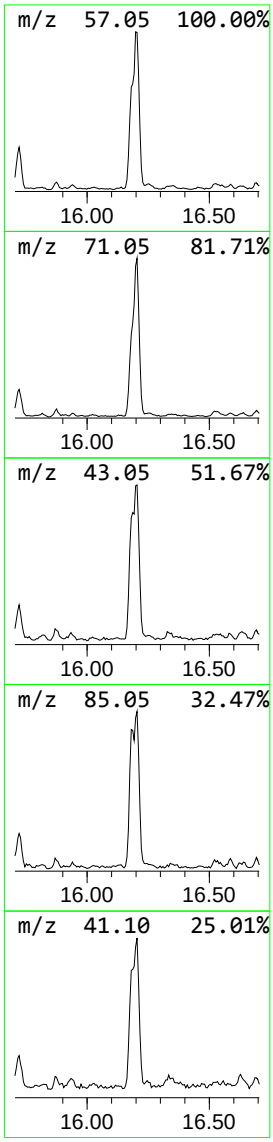
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.204	3.81 ng/ul	922324	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	93
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
3	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	86
4	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	81
5	Octane, 2-methyl-	128	C9H20	003221-61-2	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

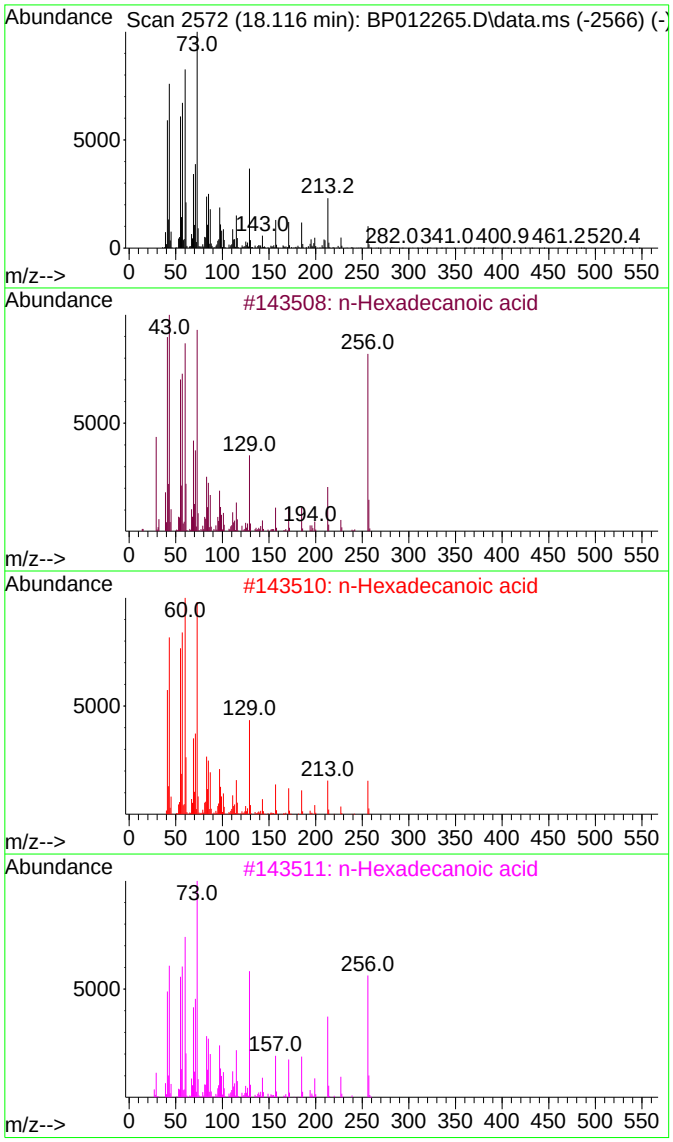
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

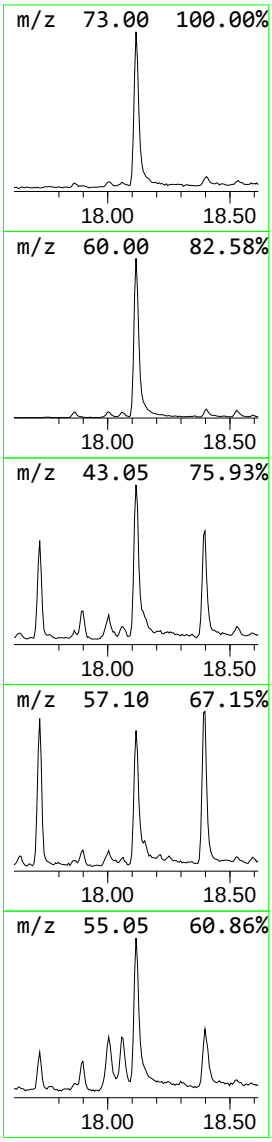
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	5.46 ng/ul	1321830	Phenanthrene-d10	17.228

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	Tridecanoic acid	214	C13H26O2	000638-53-9	96





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Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

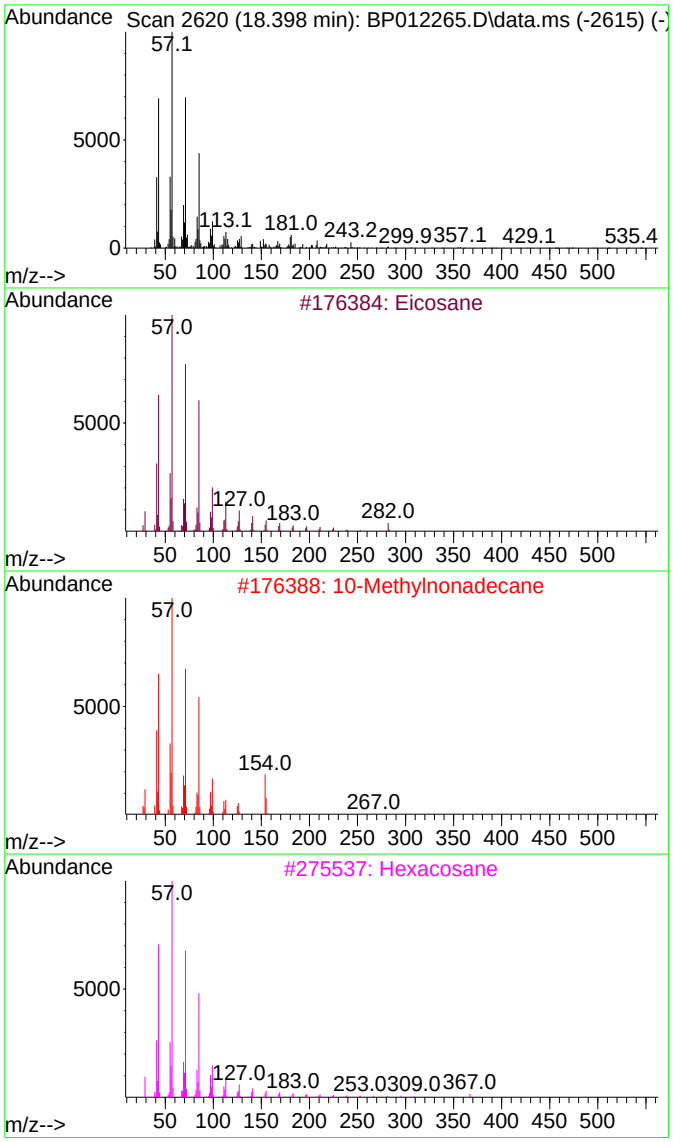
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

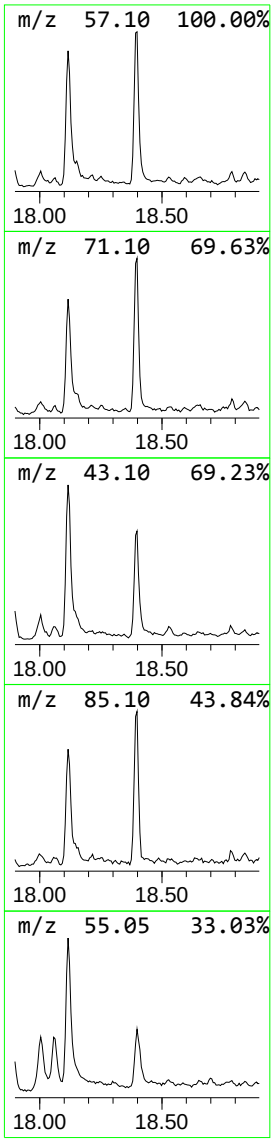
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	2.52 ng/ul	610781	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	10-Methylnonadecane	282	C20H42	056862-62-5	81
3	Hexacosane	366	C26H54	000630-01-3	81
4	Nonadecane	268	C19H40	000629-92-5	81
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	74





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Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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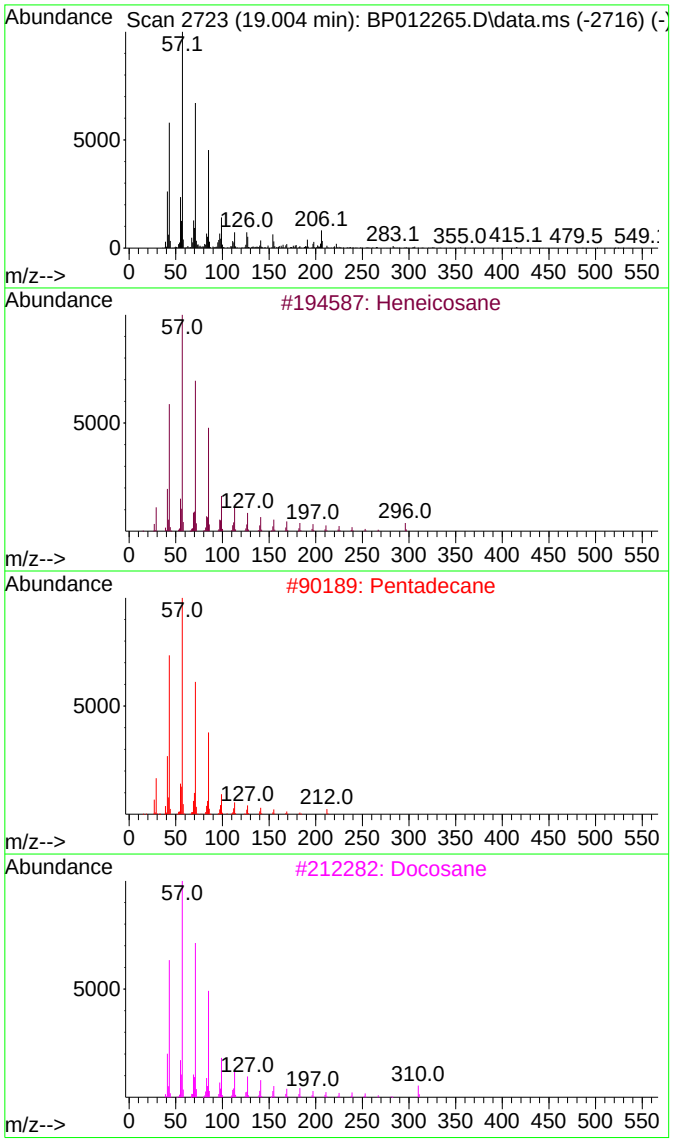
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

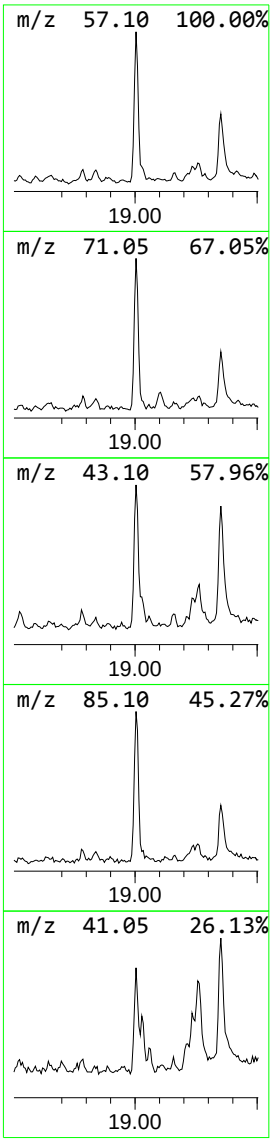
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	2.18 ng/ul	527418	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	95
2	Pentadecane	212	C15H32	000629-62-9	95
3	Docosane	310	C22H46	000629-97-0	94
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	87
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	87





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Data File : BP012265.D
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Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

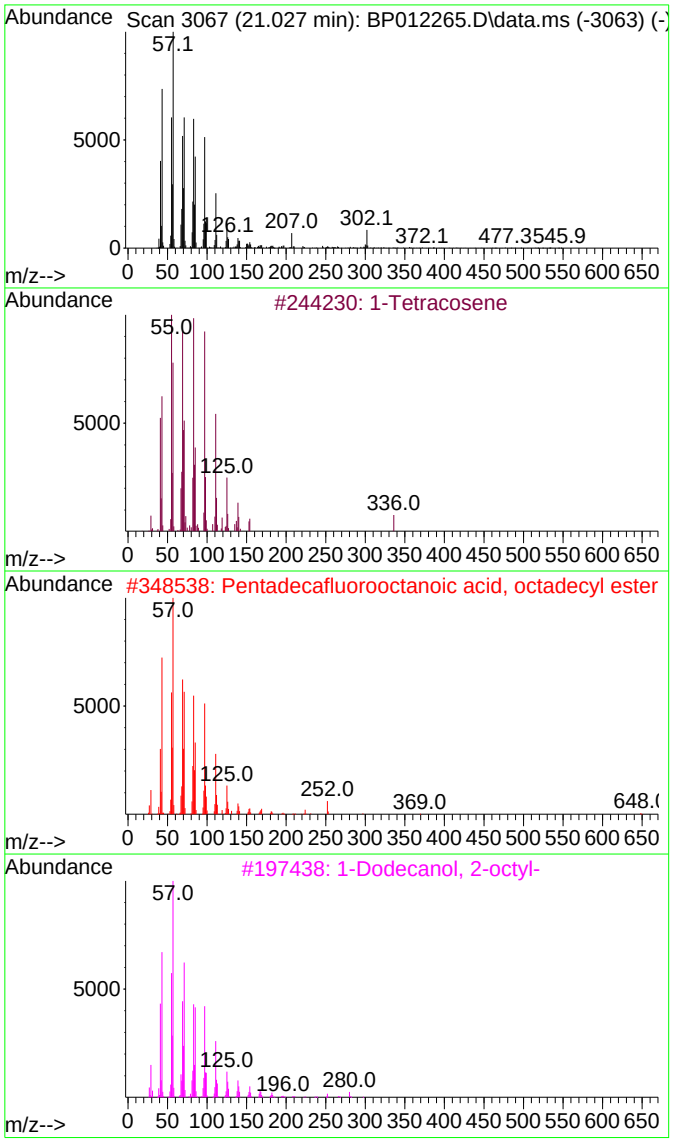
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

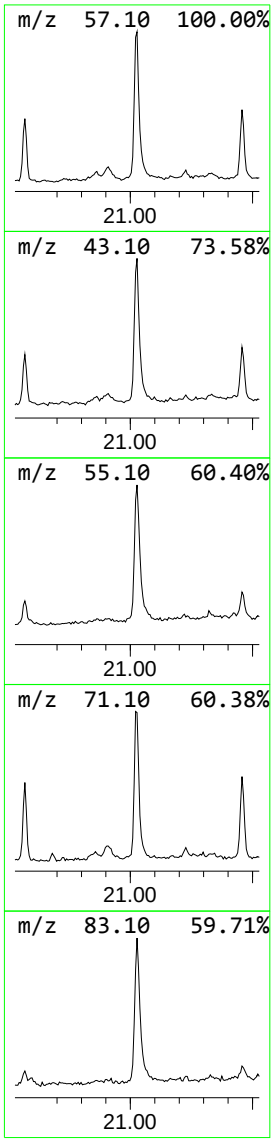
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	5.35 ng/ul	1500350	Chrysene-d12	21.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	97
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3	1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	91
4	Cyclotetradecane	196	C14H28	000295-17-0	84
5	Triacontyl acetate	480	C32H64O2	041755-58-2	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

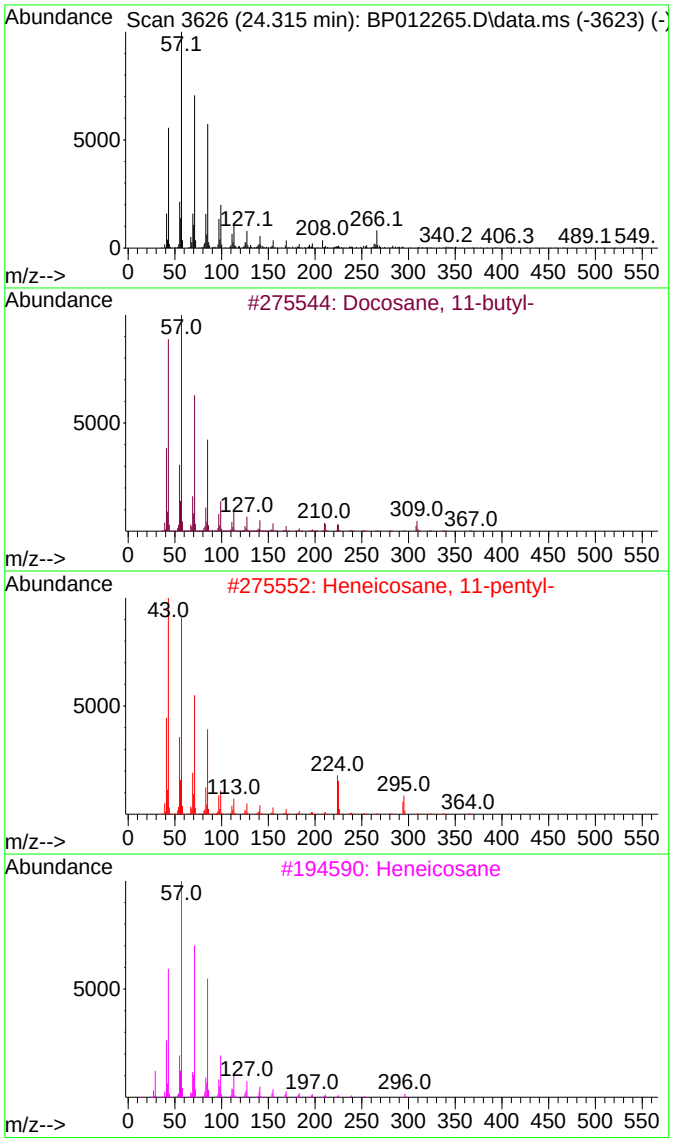
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION

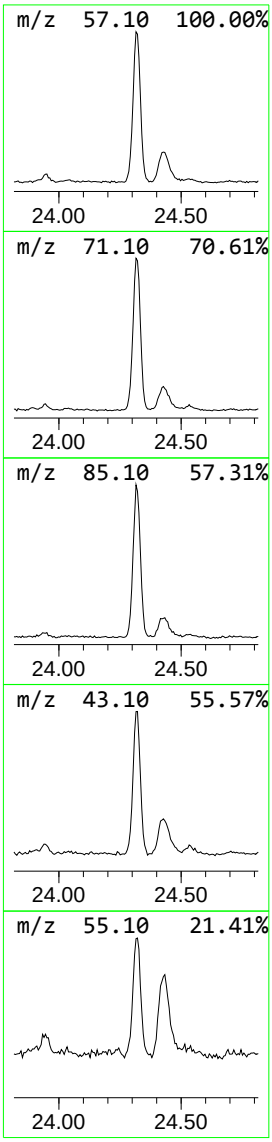
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.316	6.42 ng/ul	1311570	Perylene-d12	23.721

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90
3	Heneicosane	296	C21H44	000629-94-7	87
4	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	87
5	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012265.D
Acq On : 22 Oct 2022 02:57
Operator : CG/JU
Sample : N5064-03
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BJ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.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
3-Penten-2-ol	3.040	3.8	ng/ul	512696	1	7.816	2720380	20.0
unknown-01	3.781	4.0	ng/ul	546852	1	7.816	2720380	20.0
Ethanol, 2-(2-b...	10.404	5.1	ng/ul	1016430	2	10.622	3961830	20.0
(DEL) Alkane: S...	16.204	3.8	ng/ul	922324	4	17.228	4840870	20.0
n-Hexadecanoic ...	18.116	5.5	ng/ul	1321830	4	17.228	4840870	20.0
(DEL) Alkane: S...	18.398	2.5	ng/ul	610781	4	17.228	4840870	20.0
(DEL) Alkane: S...	19.004	2.2	ng/ul	527418	4	17.228	4840870	20.0
(DEL) Alkane: S...	21.027	5.3	ng/ul	1500350	5	21.322	5606480	20.0
(DEL) Alkane: S...	24.316	6.4	ng/ul	1311570	6	23.721	4084820	20.0