LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051822\
Data File : BM035147.D
Acq On : 18 May 2022 16:57
Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

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_M\Methods\SFAM-EPA-BM051722.M
Title : SVOA CALIBRATION

Signal : TIC: BM035147.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	26	rVB	1295784	2156036	94.14%	7.353%
2	3.375	45	49	59	rVB4	18847	43072	1.88%	0.147%
3	3.799	118	121	130	rBV2	30365	61731	2.70%	0.211%
4	3.893	133	137	145	rVB	22060	36514	1.59%	0.125%
5	4.016	155	158	164	rVB2	9887	18166	0.79%	0.062%
6	4.110	172	174	184	rVB3	13158	21720	0.95%	0.074%
7	4.669	264	269	275	rVB6	9583	18874	0.82%	0.064%
8	5.022	324	329	338	rVV	41493	64861	2.83%	0.221%
9	5.122	340	346	357	rVB9	7807	19072	0.83%	0.065%
10	6.775	622	627	632	rBV7	4115	8158	0.36%	0.028%
11	6.910	644	650	655	rVB8	3619	7690	0.34%	0.026%
12	7.075	670	678	690	rBV2	319819	633956	27.68%	2.162%
13	7.240	700	706	716	rBV	267525	431855	18.86%	1.473%
14	7.440	733	740	751	rVV	341896	563279	24.60%	1.921%
15	7.545	754	758	766	rBV4	11001	17953	0.78%	0.061%
16	7.640	772	774	781	rVB6	1620	3307	0.14%	0.011%
17	7.910	812	820	829	rBV	628799	1004775	43.87%	3.427%
18	7.981	829	832	835	rVB2	4352	5201	0.23%	0.018%
19	8.616	932	940	948	rBV	269211	440485	19.23%	1.502%
20	8.687	948	952	956	rVV3	12395	24053	1.05%	0.082%
21	8.734	956	960	966	rVB	27761	45307	1.98%	0.155%
22	9.063	1010	1016	1028	rBV	328041	558078	24.37%	1.903%
23	9.151	1028	1031	1035	rVV4	8844	13530	0.59%	0.046%
24	9.187	1035	1037	1042	rVV3	4872	7837	0.34%	0.027%
25	9.239	1042	1046	1052	rVB7	5881	11286	0.49%	0.038%
26	9.516	1090	1093	1097	rVB6	3525	5321	0.23%	0.018%
27	9.792	1132	1140	1149	rBV	225321	394290	17.22%	1.345%
28	9.857	1149	1151	1154	rVB4	3457	3524	0.15%	0.012%
29	9.939	1161	1165	1167	rBV4	3167	4243	0.19%	0.014%
30	10.157	1197	1202	1209	rVB8	6535	10551	0.46%	0.036%
31	10.251	1216	1218	1222	rBV4	2343	3483	0.15%	0.012%
32	10.328	1224	1231	1243	rVV	322728	568110	24.81%	1.938%
33	10.628	1277	1282	1287	rVB5	5422	9151	0.40%	0.031%
34	10.710	1289	1296	1307	rBV	799672	1338794	58.46%	4.566%
35	10.839	1310	1318	1334	rBV	297166	576738	25.18%	1.967%
36	11.063	1352	1356	1363	rBV7	3644	7849	0.34%	0.027%

LSC Area Percent Report

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Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Title : SVOA CALIBRATION

37	11.263	1383	1390	1396	rVB8	2615	7656	0.33%	0.026%
38	11.422	1415	1417	1420	rVB4	3465	3345	0.15%	0.011%
39	11.510	1426	1432	1436	rBV6	3935	8568	0.37%	0.029%
40	11.751	1469	1473	1483	rVB5	7302	14419	0.63%	0.049%
41	12.057	1520	1525	1531	rVB	15306	24000	1.05%	0.082%
42	12.145	1534	1540	1547	rVB2	14929	23348	1.02%	0.080%
43	12.298	1560	1566	1571	rVB4	7836	14863	0.65%	0.051%
44	12.375	1574	1579	1582	rBV5	3484	5981	0.26%	0.020%
45	12.775	1641	1647	1650	rBV7	3059	6038	0.26%	0.021%
46	12.828	1650	1656	1658	rVV6	3527	5871	0.26%	0.020%
47	12.886	1661	1666	1668	rBV6	2619	3583	0.16%	0.012%
48	12.951	1676	1677	1683	rVB5	3619	4714	0.21%	0.016%
49	13.045	1690	1693	1698	rBV6	2438	4317	0.19%	0.015%
50	13.098	1698	1702	1707	rVB	13160	19585	0.86%	0.067%
51	13.157	1709	1712	1718	rBV7	4182	6651	0.29%	0.023%
52	13.380	1743	1750	1756	rBV	35437	56262	2.46%	0.192%
53	13.475	1763	1766	1771	rVB5	3924	4451	0.19%	0.015%
54	13.551	1776	1779	1783	rVB4	3766	4781	0.21%	0.016%
55	13.663	1793	1798	1802	rVB7	3240	4474	0.20%	0.015%
56	13.704	1802	1805	1807	rBV4	3375	3988	0.17%	0.014%
57	13.822	1820	1825	1829	rBV8	4012	6097	0.27%	0.021%
58	13.869	1829	1833	1838	rBV6	9701	18697	0.82%	0.064%
59	13.951	1841	1847	1855	rVV	722803	963316	42.06%	3.285%
60	14.039	1859	1862	1866	rVB2	28180	35359	1.54%	0.121%
61	14.080	1866	1869	1875	rVB5	9799	15637	0.68%	0.053%
62	14.157	1878	1882	1885	rVB3	5646	7665	0.33%	0.026%
63	14.233	1885	1895	1904	rBV	777320	1183079	51.66%	4.035%
64	14.322	1908	1910	1913	rVB3	6142	5552	0.24%	0.019%
65	14.380	1916	1920	1927	rVB9	11963	22176	0.97%	0.076%
66	14.457	1927	1933	1938	rBV	82598	115305	5.03%	0.393%
67	14.545	1940	1948	1954	rBV	1148120	1699209	74.20%	5.795%
68	14.769	1982	1986	1988	rBV5	5716	6440	0.28%	0.022%
69	14.898	2006	2008	2009	rBV2	6127	5247	0.23%	0.018%
70	14.957	2014	2018	2023	rVB3	20625	32073	1.40%	0.109%
71	15.010	2025	2027	2032	rVB5	9677	11211	0.49%	0.038%
72	15.074	2034	2038	2043	rVB2	37134	50876	2.22%	0.174%
73	15.139	2046	2049	2052	rBV2	12562	17027	0.74%	0.058%
74	15.204	2058	2060	2065	rVB6	11249	14258	0.62%	0.049%
75	15.263	2068	2070	2076	rVV7	7452	10716	0.47%	0.037%
76	15.339	2080	2083	2089	rVB6	24123	35499	1.55%	0.121%

LSC Area Percent Report

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Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

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
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Title : SVOA CALIBRATION

77	15.404	2089	2094	2098	rBV	154625	183299	8.00%	0.625%
78	15.533	2110	2116	2123	rVB	1009247	1462256	63.85%	4.987%
79	15.810	2154	2163	2169	rBV	121335	235058	10.26%	0.802%
80	15.963	2185	2189	2194	rBV3	33914	66614	2.91%	0.227%
81	16.027	2196	2200	2203	rVV3	49084	69188	3.02%	0.236%
82	16.192	2225	2228	2230	rBV4	15243	21869	0.95%	0.075%
83	16.268	2236	2241	2243	rBV	282530	405214	17.69%	1.382%
84	16.292	2243	2245	2250	rVB	226893	260265	11.36%	0.888%
85	17.057	2371	2375	2379	rBV	258688	305731	13.35%	1.043%
86	17.115	2380	2385	2392	rVB	436658	649869	28.38%	2.216%
87	17.286	2409	2414	2420	rBV	1272563	1814420	79.23%	6.188%
88	17.386	2425	2431	2438	rVV	1036972	1488524	65.00%	5.077%
89	17.721	2485	2488	2492	rVB	100980	130028	5.68%	0.443%
90	17.798	2498	2501	2506	rVB	285192	333698	14.57%	1.138%
91	18.192	2564	2568	2574	rBV2	218750	338917	14.80%	1.156%
92	18.492	2616	2619	2624	rBV	206843	250771	10.95%	0.855%
93	19.127	2724	2727	2731	rBV	158179	178691	7.80%	0.609%
94	19.468	2782	2785	2788	rBV2	164819	213854	9.34%	0.729%
95	19.680	2816	2821	2828	rBV	1178796	1780904	77.76%	6.074%
96	21.468	3121	3125	3136	rVB	1511477	1942334	84.81%	6.624%
97	23.703	3501	3505	3513	rVB2	719923	1277738	55.79%	4.358%
98	23.862	3527	3532	3542	rVB	1134379	2290193	100.00%	7.811%

Sum of corrected areas: 29320619

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Instrument :
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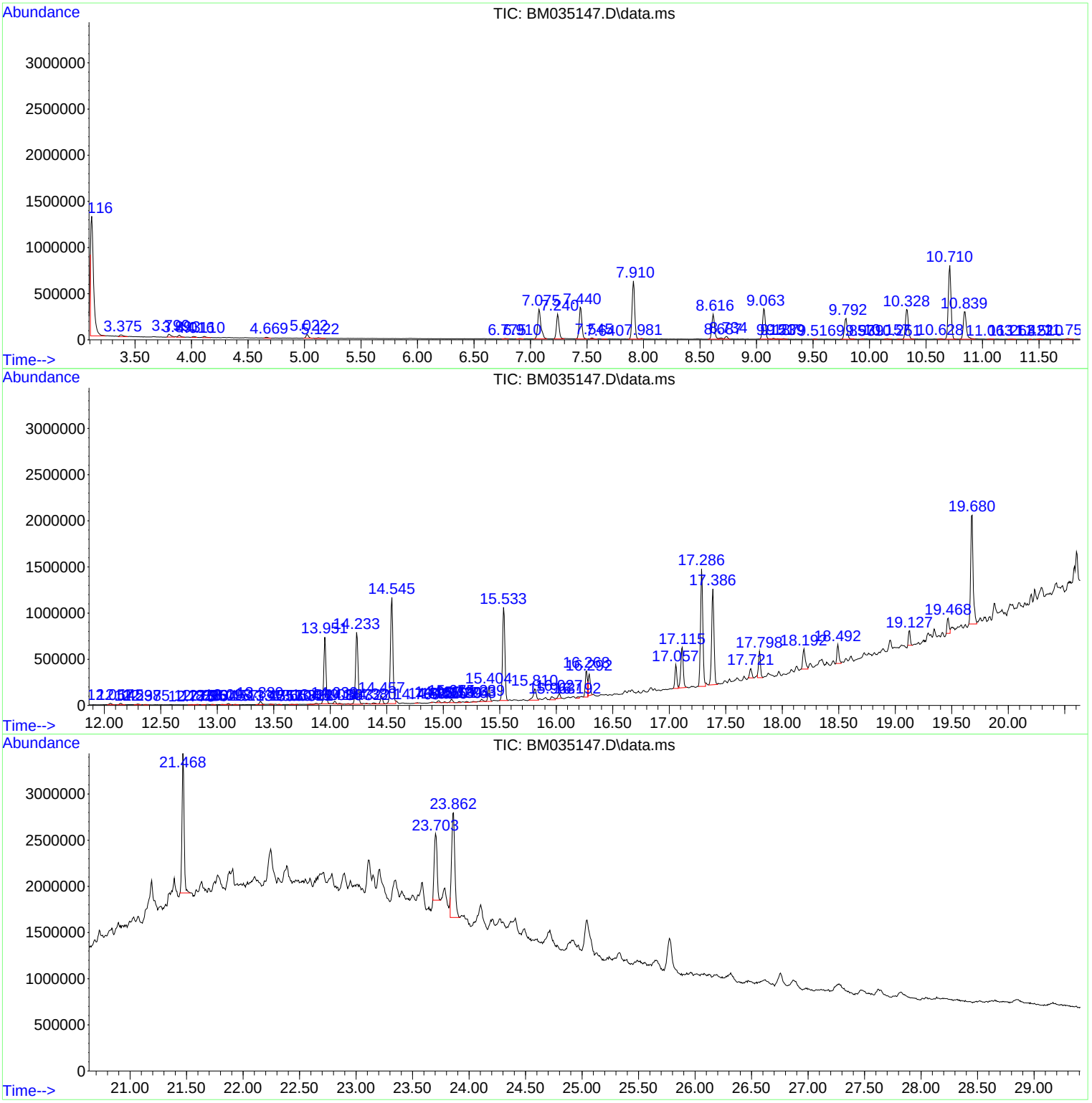
ClientSampleId :
EW4G7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

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



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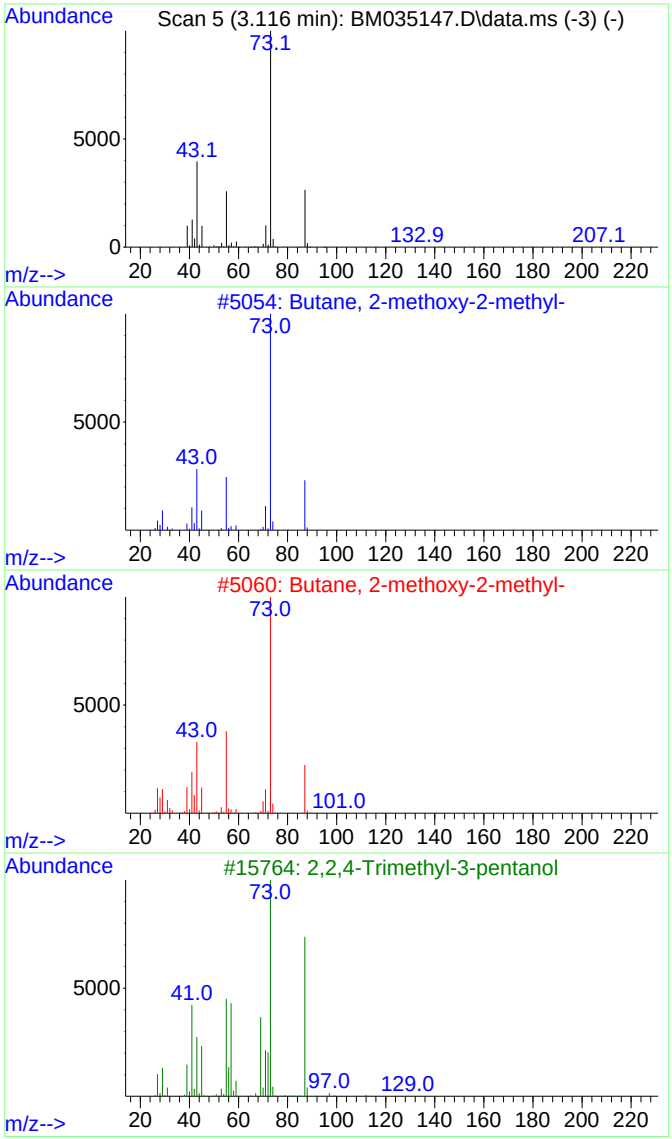
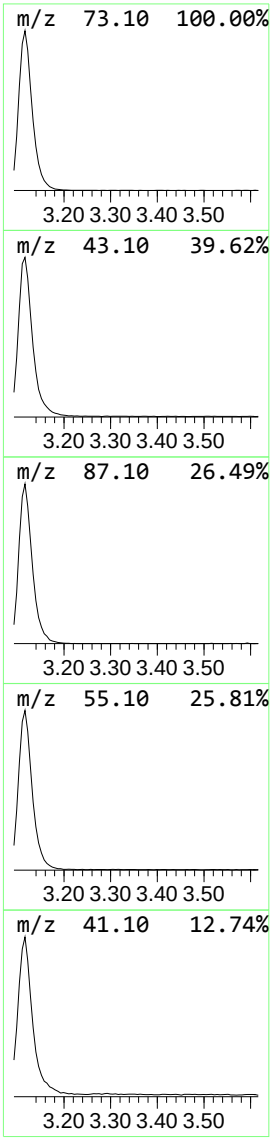
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.116	42.92 ng/ul	2156040	1,4-Dichlorobenzene-d4	7.910

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	50
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12





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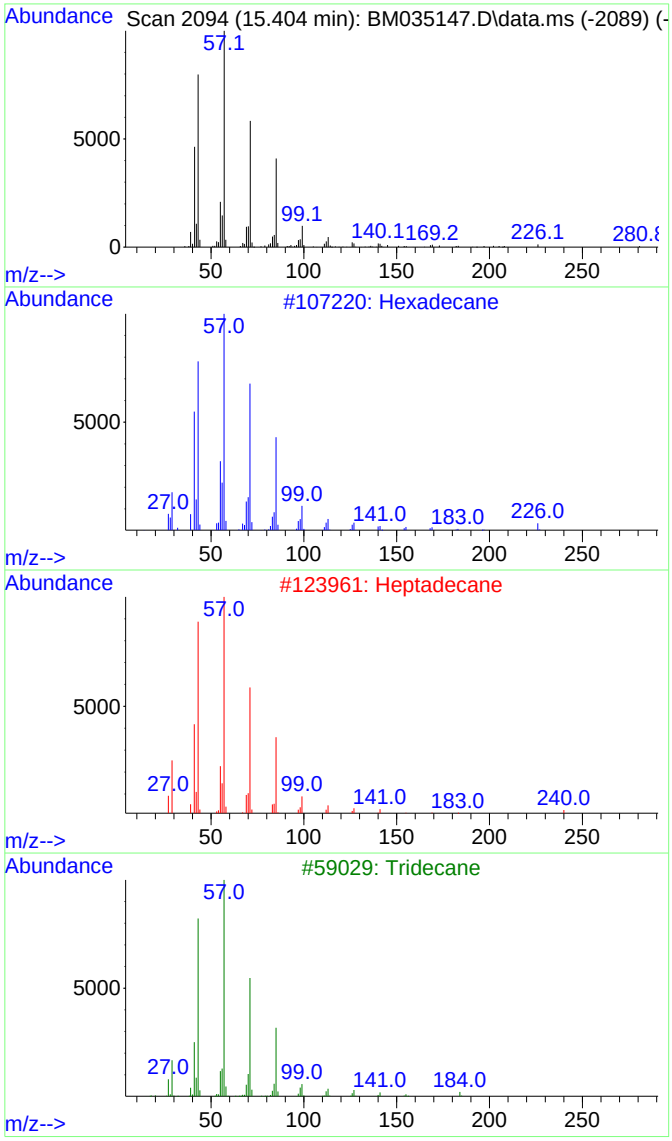
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

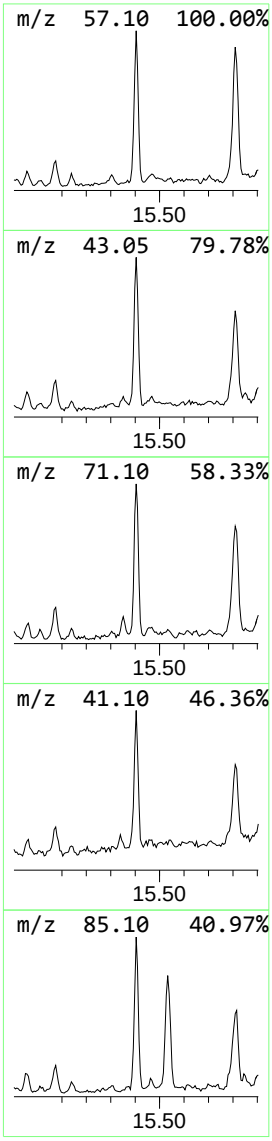
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	2.16 ng/ul	183299	Acenaphthene-d10	14.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Heptadecane	240	C17H36	000629-78-7	90
3	Tridecane	184	C13H28	000629-50-5	86
4	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1010309-12-5	80
5	Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	80





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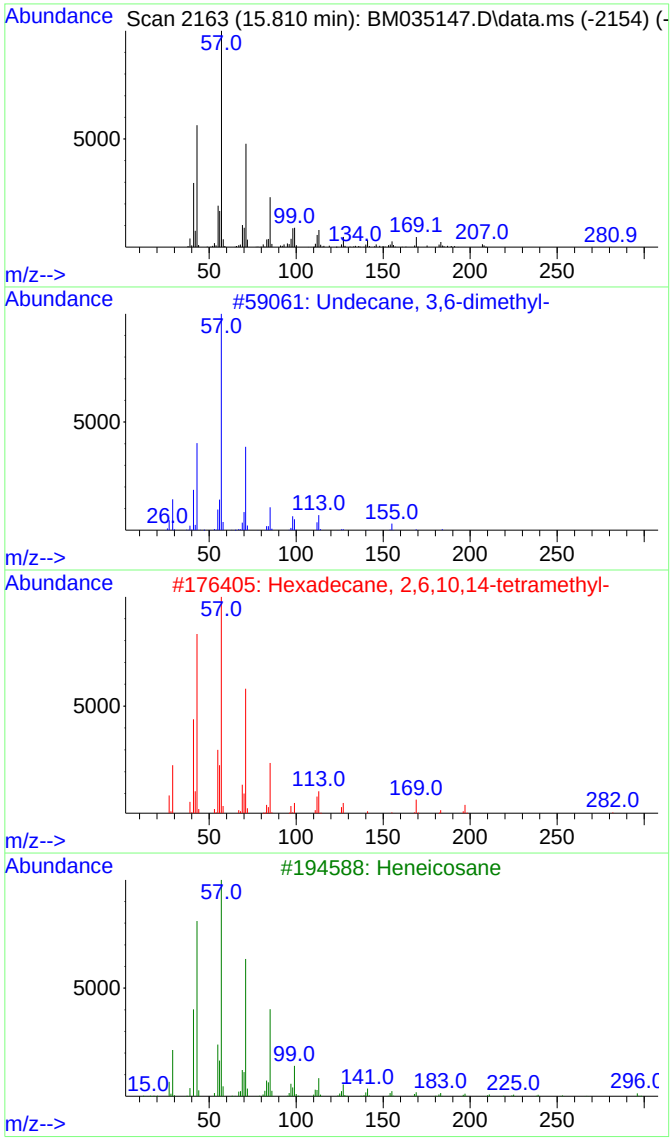
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

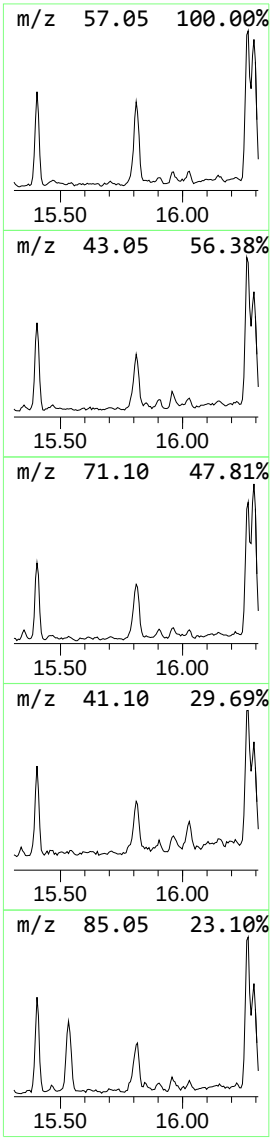
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.810	2.77 ng/ul	235058	Acenaphthene-d10	14.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	81
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80
3	Heneicosane	296	C21H44	000629-94-7	72
4	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	72
5	Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	68





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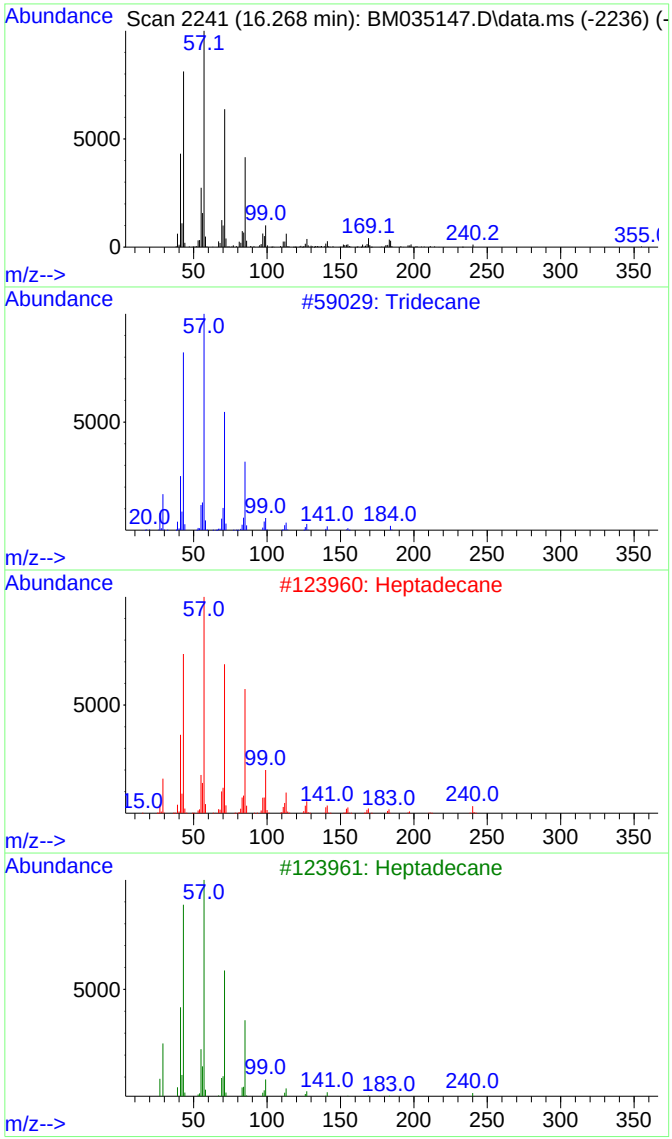
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

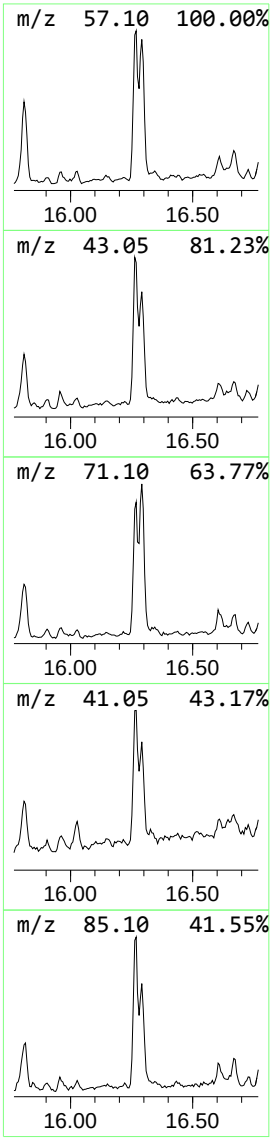
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.269	4.47 ng/ul	405214	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	96
2	Heptadecane	240	C17H36	000629-78-7	95
3	Heptadecane	240	C17H36	000629-78-7	95
4	Heptadecane	240	C17H36	000629-78-7	93
5	Nonadecane	268	C19H40	000629-92-5	87





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Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

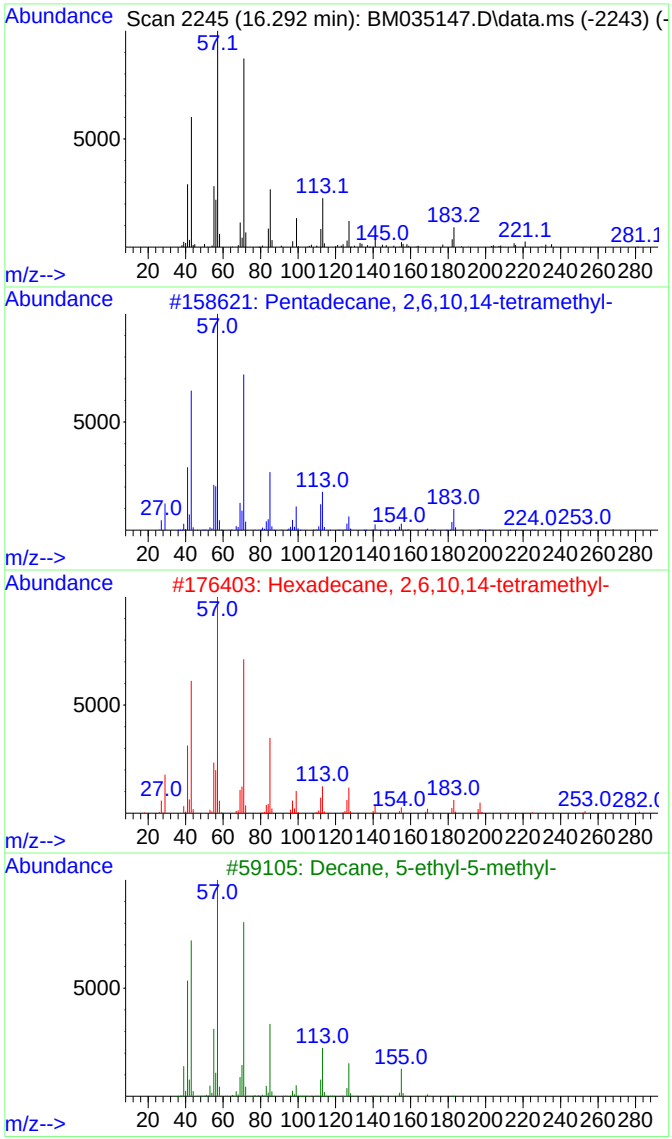
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

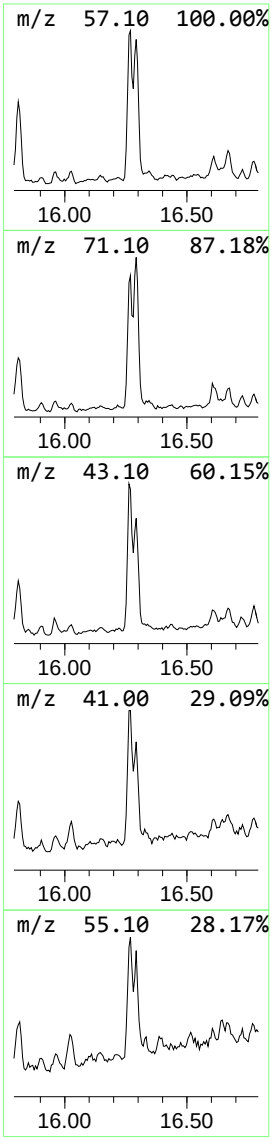
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.292	2.87 ng/ul	260265	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
3	Decane, 5-ethyl-5-methyl-	184	C13H28	017312-74-2	64
4	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
5	Pentadecane, 2-methyl-	226	C16H34	001560-93-6	59





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Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

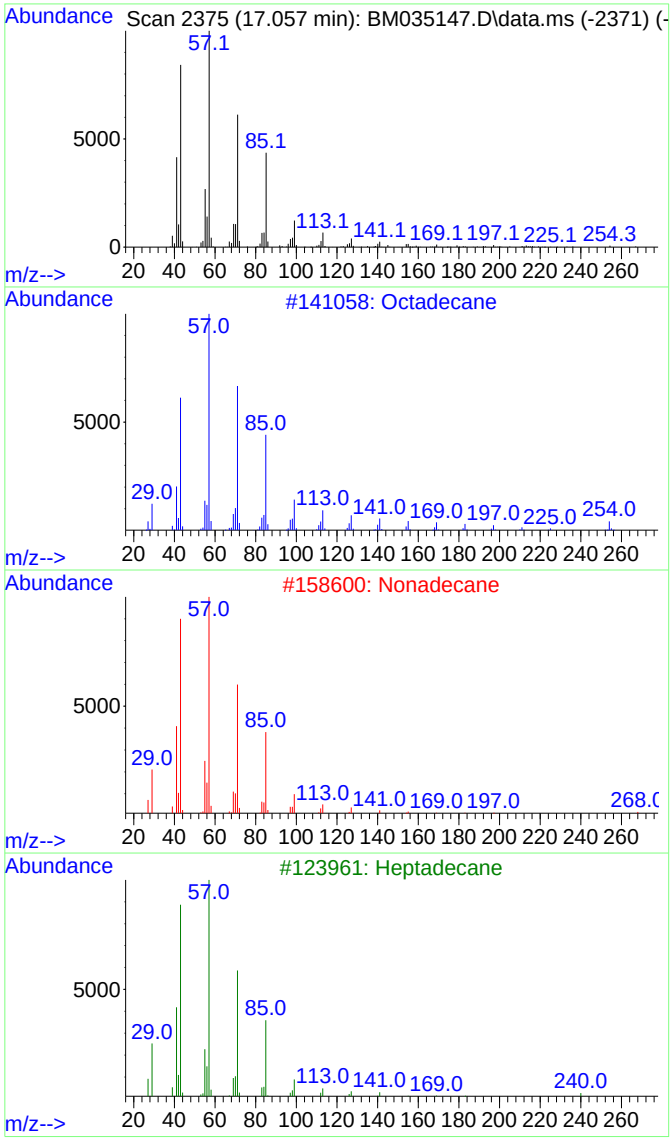
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

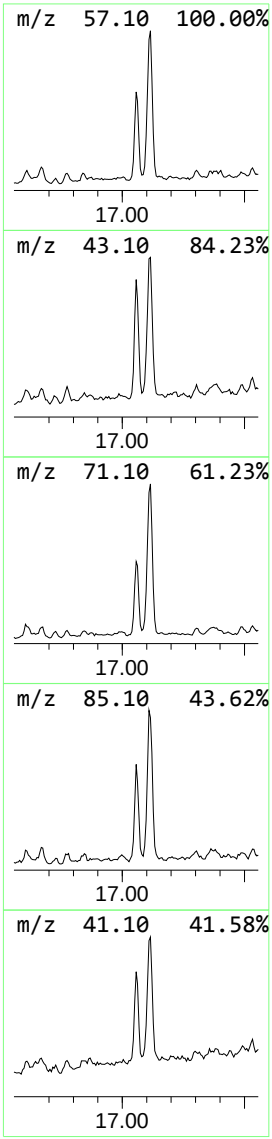
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.057	3.37 ng/ul	305731	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	Nonadecane	268	C19H40	000629-92-5	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
5	Heptacosane	380	C27H56	000593-49-7	90





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Acq On : 18 May 2022 16:57
Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

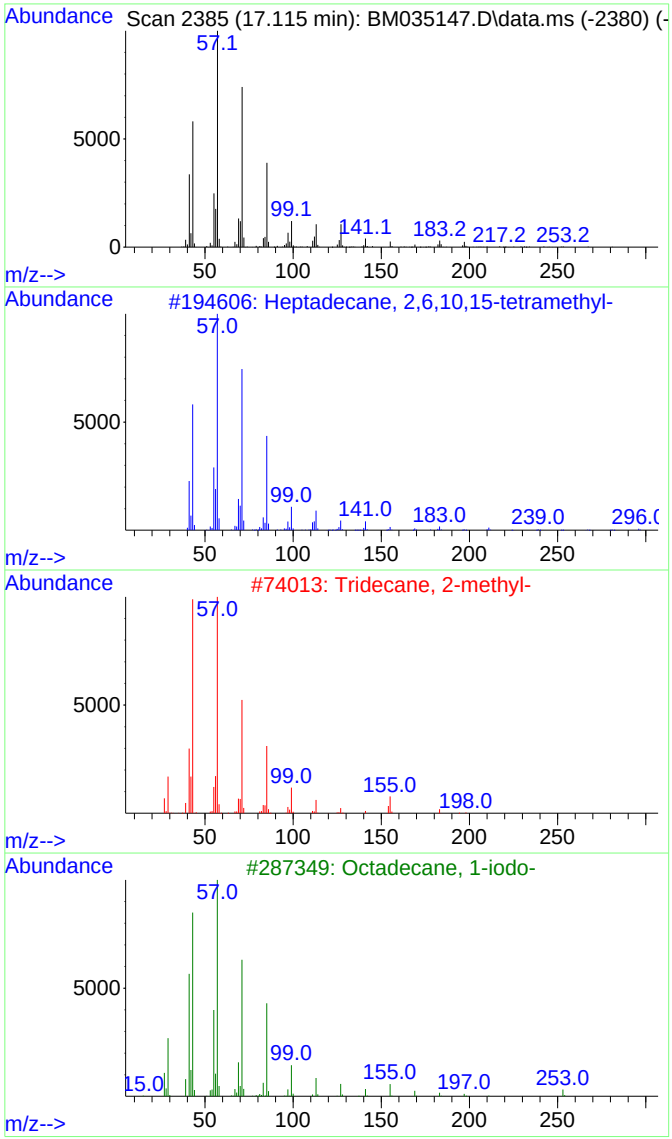
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

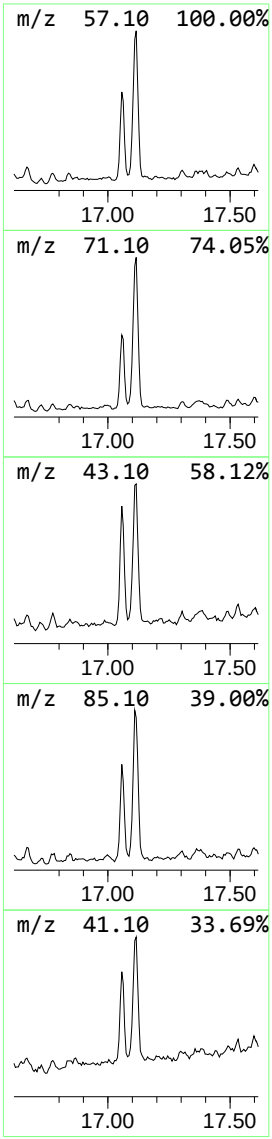
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.115	7.16 ng/ul	649869	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2	Tridecane, 2-methyl-	198	C14H30	001560-96-9	90
3	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90
4	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
5	Tridecane, 4-methyl-	198	C14H30	026730-12-1	87





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Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

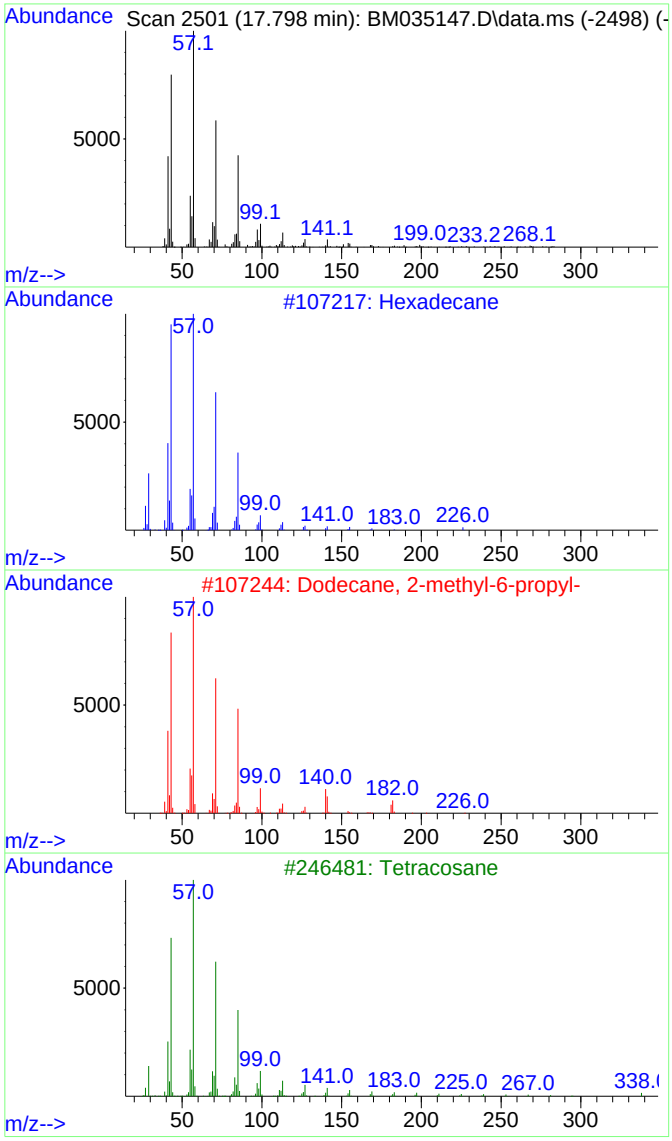
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

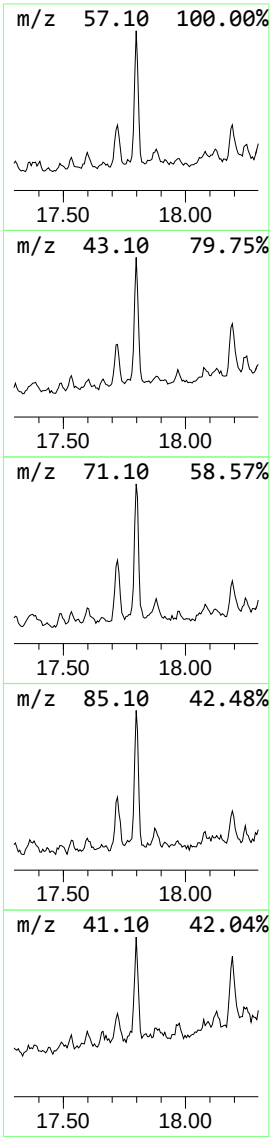
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.798	3.68 ng/ul	333698	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
3	Tetracosane	338	C24H50	000646-31-1	91
4	Decane, 1-iodo-	268	C10H21I	002050-77-3	90
5	Hexacosane	366	C26H54	000630-01-3	90





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Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

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

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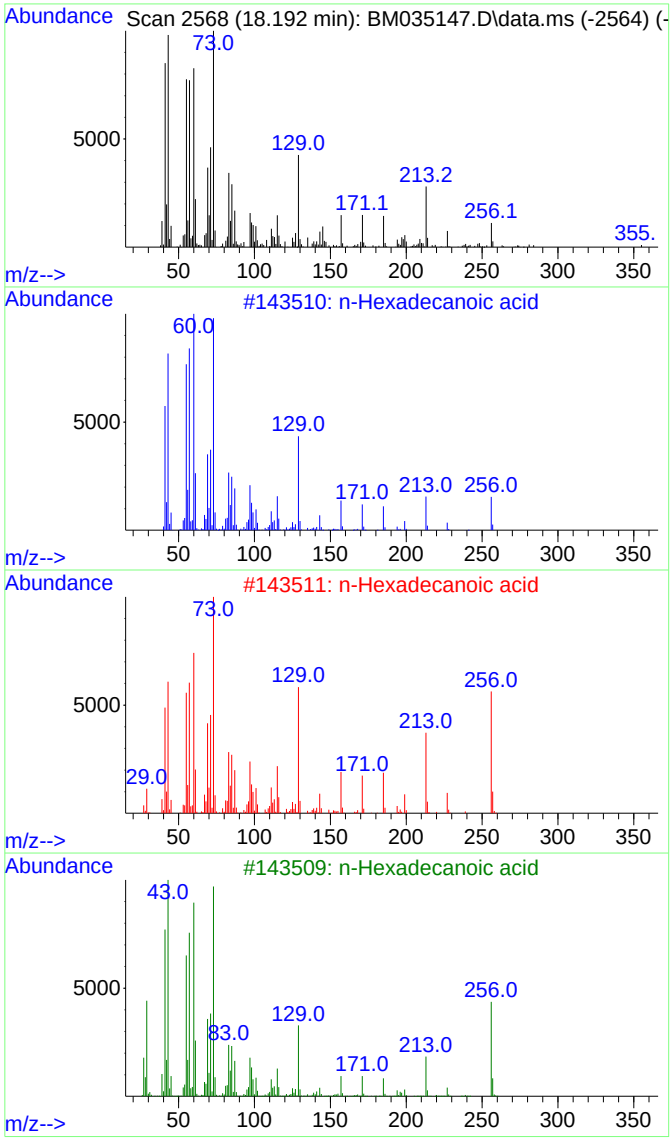
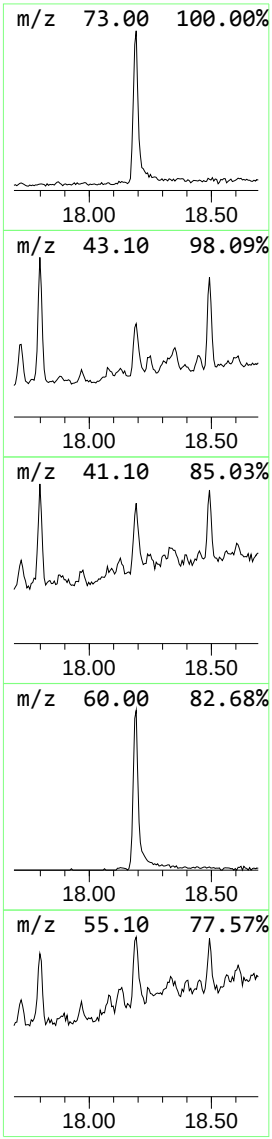
Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	3.74 ng/ul	338917	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91





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Data File : BM035147.D
Acq On : 18 May 2022 16:57
Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

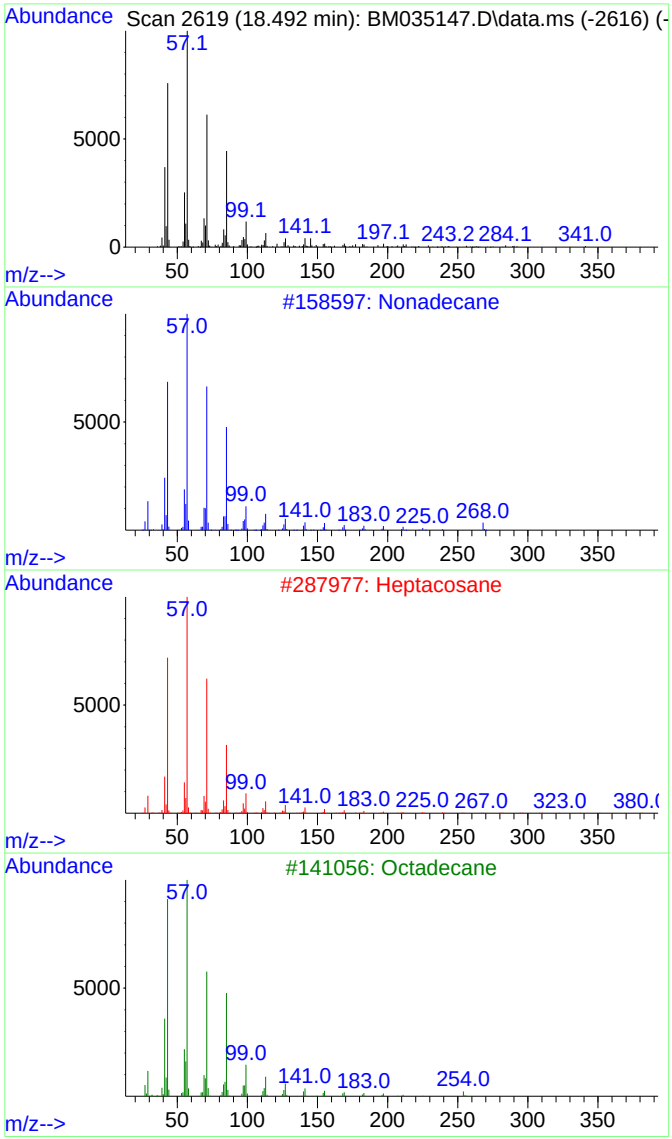
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

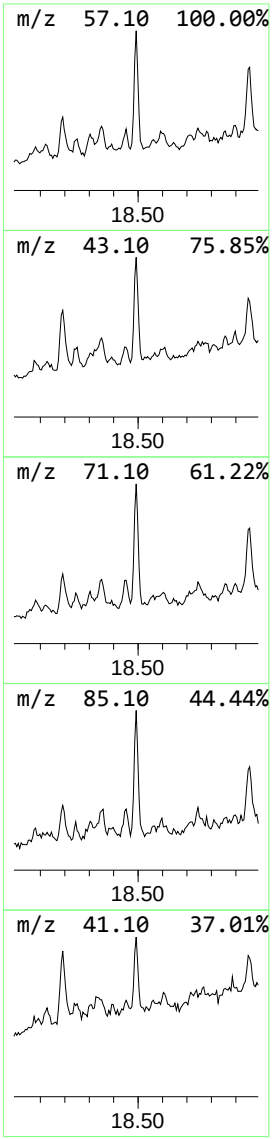
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.492	2.76 ng/ul	250771	Phenanthrene-d10	17.286

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	91
2	Heptacosane	380	C27H56	000593-49-7	91
3	Octadecane	254	C18H38	000593-45-3	87
4	Eicosane	282	C20H42	000112-95-8	87
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	87





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Data File : BM035147.D
Acq On : 18 May 2022 16:57
Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
Quant Title : SVOA CALIBRATION

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

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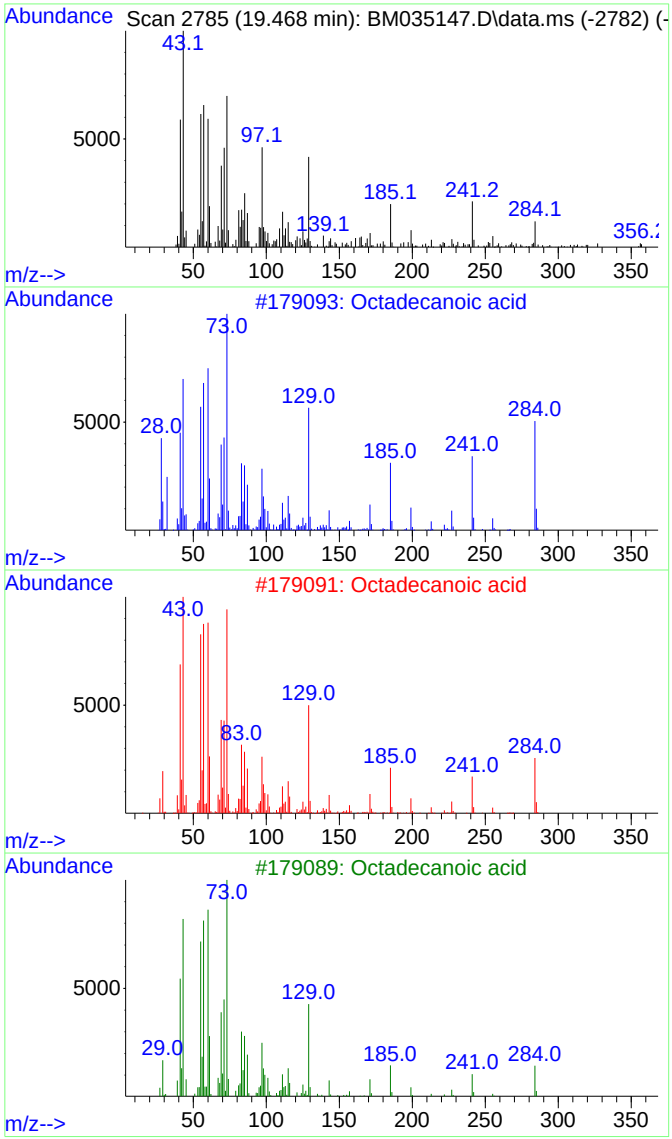
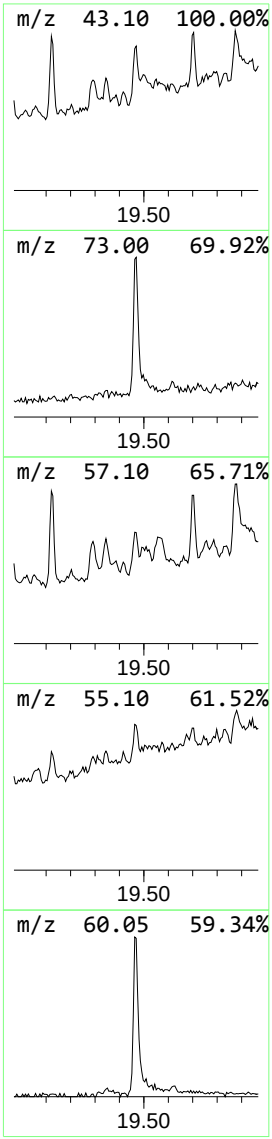
Peak Number 11 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.468	2.20 ng/ul	213854	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	92
2	Octadecanoic acid	284	C18H36O2	000057-11-4	91
3	Octadecanoic acid	284	C18H36O2	000057-11-4	89
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	60
5	Octadecanoic acid	284	C18H36O2	000057-11-4	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051822\
Data File : BM035147.D
Acq On : 18 May 2022 16:57
Operator : CG/JU
Sample : N2766-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4G7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.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
Butane, 2-metho...	3.116	42.9	ng/ul	2156040	1	7.910	1004780	20.0
(DEL) Alkane: S...	15.404	2.2	ng/ul	183299	3	14.545	1699210	20.0
(DEL) Alkane: S...	15.810	2.8	ng/ul	235058	3	14.545	1699210	20.0
(DEL) Alkane: S...	16.269	4.5	ng/ul	405214	4	17.286	1814420	20.0
(DEL) Alkane: S...	16.292	2.9	ng/ul	260265	4	17.286	1814420	20.0
(DEL) Alkane: S...	17.057	3.4	ng/ul	305731	4	17.286	1814420	20.0
(DEL) Alkane: S...	17.115	7.2	ng/ul	649869	4	17.286	1814420	20.0
(DEL) Alkane: S...	17.798	3.7	ng/ul	333698	4	17.286	1814420	20.0
n-Hexadecanoic ...	18.192	3.7	ng/ul	338917	4	17.286	1814420	20.0
(DEL) Alkane: S...	18.492	2.8	ng/ul	250771	4	17.286	1814420	20.0
Octadecanoic acid	19.468	2.2	ng/ul	213854	5	21.468	1942330	20.0