LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015648.D
Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCFK8

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-BP060723.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP015648.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	33	36	43	rVB	22824	29634	0.84%	0.076%
2	3.846	109	112	124	rVV	184808	303782	8.64%	0.780%
3	3.946	125	129	137	rVV	55279	87327	2.48%	0.224%
4	4.069	146	150	158	rVB3	18436	30870	0.88%	0.079%
5	4.169	164	167	173	rVB3	10829	14645	0.42%	0.038%
6	5.240	346	349	358	rVB9	7054	14848	0.42%	0.038%
7	6.046	481	486	490	rVB4	9635	13306	0.38%	0.034%
8	6.299	523	529	540	rBV3	11718	36854	1.05%	0.095%
9	6.705	596	598	603	rVV5	8277	14215	0.40%	0.036%
10	7.240	683	689	705	rBV	325167	621234	17.67%	1.594%
11	7.399	710	716	731	rVB	276513	475848	13.54%	1.221%
12	7.605	745	751	760	rBV	388690	657434	18.70%	1.687%
13	7.675	761	763	767	rVV2	15929	20122	0.57%	0.052%
14	8.081	825	832	845	rVV	509041	878659	24.99%	2.255%
15	8.728	934	942	946	rBV10	6456	15394	0.44%	0.040%
16	8.804	946	955	970	rBV	370158	749661	21.33%	1.924%
17	9.257	1023	1032	1046	rBV	364342	696754	19.82%	1.788%
18	9.410	1054	1058	1067	rVB	7582	20617	0.59%	0.053%
19	9.987	1148	1156	1167	rBV	304847	573837	16.32%	1.473%
20	10.534	1242	1249	1270	rBV	355348	841440	23.94%	2.159%
21	10.857	1298	1304	1306	rBV3	13982	21811	0.62%	0.056%
22	10.904	1306	1312	1319	rVV	737171	1236268	35.17%	3.172%
23	10.957	1319	1321	1327	rVV	39071	53287	1.52%	0.137%
24	11.063	1331	1339	1358	rVV	110559	315586	8.98%	0.810%
25	11.904	1476	1482	1486	rBV3	20483	31799	0.90%	0.082%
26	12.304	1543	1550	1556	rBV3	16948	34806	0.99%	0.089%
27	12.493	1579	1582	1589	rVB8	8085	16017	0.46%	0.041%
28	12.569	1589	1595	1603	rBV2	23945	45134	1.28%	0.116%
29	12.787	1627	1632	1636	rVV	21060	33827	0.96%	0.087%
30	13.104	1680	1686	1690	rVV8	5408	12986	0.37%	0.033%
31	13.240	1703	1709	1715	rBV	18279	27008	0.77%	0.069%
32	13.528	1753	1758	1763	rBV2	26595	41975	1.19%	0.108%
33	13.604	1767	1771	1779	rVB2	21332	41171	1.17%	0.106%
34	13.892	1812	1820	1821	rBV8	10690	16696	0.47%	0.043%
35	14.045	1840	1846	1851	rBV4	33561	65506	1.86%	0.168%
36	14.134	1851	1861	1867	rBV	836610	1245039	35.42%	3.195%

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Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

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-BP060723.MA.M
Title : SVOA CALIBRATION

37	14.181	1867	1869	1873	rVB2	31873	37124	1.06%	0.095%
38	14.287	1883	1887	1889	rBV3	10663	12846	0.37%	0.033%
39	14.422	1903	1910	1924	rBV2	932981	1495844	42.55%	3.838%
40	14.545	1926	1931	1935	rVV7	12061	18093	0.51%	0.046%
41	14.598	1935	1940	1945	rVB	30731	39301	1.12%	0.101%
42	14.675	1949	1953	1955	rBV4	14009	19125	0.54%	0.049%
43	14.728	1955	1962	1969	rVV	975848	1477664	42.03%	3.792%
44	14.786	1969	1972	1982	rVB2	32927	60294	1.72%	0.155%
45	14.969	1990	2003	2024	rBV2	99812	391771	11.14%	1.005%
46	15.134	2026	2031	2037	rVV3	36444	76623	2.18%	0.197%
47	15.198	2040	2042	2048	rVV4	12565	20929	0.60%	0.054%
48	15.345	2062	2067	2070	rBV3	13490	22558	0.64%	0.058%
49	15.392	2073	2075	2080	rVV2	9990	12987	0.37%	0.033%
50	15.504	2087	2094	2098	rBV9	23125	44509	1.27%	0.114%
51	15.551	2098	2102	2106	rVB	36846	46943	1.34%	0.120%
52	15.722	2124	2131	2137	rBV	1067950	1581953	45.00%	4.059%
53	15.863	2150	2155	2165	rBV	158869	313211	8.91%	0.804%
54	16.086	2186	2193	2204	rBV	370813	557616	15.86%	1.431%
55	16.222	2213	2216	2224	rVB2	16078	28419	0.81%	0.073%
56	16.433	2245	2252	2267	rVB2	62182	173574	4.94%	0.445%
57	16.575	2273	2276	2281	rBV5	15750	24726	0.70%	0.063%
58	16.651	2284	2289	2294	rVB	41290	60656	1.73%	0.156%
59	17.016	2345	2351	2354	rBV5	17589	35585	1.01%	0.091%
60	17.051	2355	2357	2361	rVB4	11609	14529	0.41%	0.037%
61	17.145	2368	2373	2380	rBV2	37569	63915	1.82%	0.164%
62	17.216	2380	2385	2390	rBV	52093	79263	2.25%	0.203%
63	17.269	2390	2394	2397	rVB6	16391	21638	0.62%	0.056%
64	17.310	2397	2401	2406	rVB	23225	32944	0.94%	0.085%
65	17.363	2406	2410	2418	rBV4	41784	70726	2.01%	0.181%
66	17.433	2418	2422	2425	rBV5	24877	33416	0.95%	0.086%
67	17.486	2425	2431	2435	rBV	951977	1344379	38.24%	3.450%
68	17.528	2435	2438	2443	rVB	398768	520767	14.81%	1.336%
69	17.586	2443	2448	2453	rBV	894591	1274974	36.27%	3.272%
70	17.898	2497	2501	2507	rBV2	41993	73245	2.08%	0.188%
71	17.951	2507	2510	2514	rVB	49597	53404	1.52%	0.137%
72	18.233	2555	2558	2563	rVB6	26524	34164	0.97%	0.088%
73	18.292	2564	2568	2573	rBV	137822	203775	5.80%	0.523%
74	18.345	2573	2577	2583	rVV2	517159	738558	21.01%	1.895%
75	18.398	2583	2586	2596	rVB2	104146	184636	5.25%	0.474%
76	18.539	2605	2610	2613	rBV3	73287	125010	3.56%	0.321%

LSC Area Percent Report

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Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Title : SVOA CALIBRATION

77	18.622	2620	2624	2632	rBV5	54690	113959	3.24%	0.292%
78	18.827	2656	2659	2663	rBV	53329	68127	1.94%	0.175%
79	18.880	2665	2668	2673	rVB3	52325	78399	2.23%	0.201%
80	19.227	2723	2727	2732	rVB2	111415	130208	3.70%	0.334%
81	19.374	2750	2752	2758	rVB2	54720	84091	2.39%	0.216%
82	19.439	2760	2763	2764	rBV2	35749	41206	1.17%	0.106%
83	19.486	2767	2771	2777	rVB	1004191	1306283	37.16%	3.352%
84	19.574	2783	2786	2793	rBV4	95717	148610	4.23%	0.381%
85	19.816	2819	2827	2829	rBV	1149790	1691732	48.12%	4.341%
86	19.839	2829	2831	2839	rVB	720539	909763	25.88%	2.335%
87	19.910	2840	2843	2850	rBV3	49289	110492	3.14%	0.284%
88	20.298	2907	2909	2918	rVB2	125472	247306	7.04%	0.635%
89	20.786	2989	2992	2996	rVB2	106250	111628	3.18%	0.286%
90	21.057	3035	3038	3044	rBV	143705	191344	5.44%	0.491%
91	21.251	3067	3071	3076	rVB2	295588	430184	12.24%	1.104%
92	21.451	3102	3105	3111	rVB	315979	347390	9.88%	0.891%
93	21.574	3120	3126	3130	rBV2	1105035	1926139	54.79%	4.943%
94	21.616	3130	3133	3140	rVB	429126	552171	15.71%	1.417%
95	23.286	3414	3417	3422	rVB	355510	528405	15.03%	1.356%
96	23.345	3422	3427	3442	rVB	499641	1697269	48.28%	4.355%
97	23.962	3527	3532	3537	rBV2	887205	1579587	44.93%	4.053%
98	24.127	3555	3560	3567	rBV	653983	1308730	37.23%	3.358%
99	24.633	3639	3646	3654	rVB2	1582819	3515352	100.00%	9.021%
100	24.745	3660	3665	3677	rVB	482036	1094098	31.12%	2.808%

Sum of corrected areas: 38969564

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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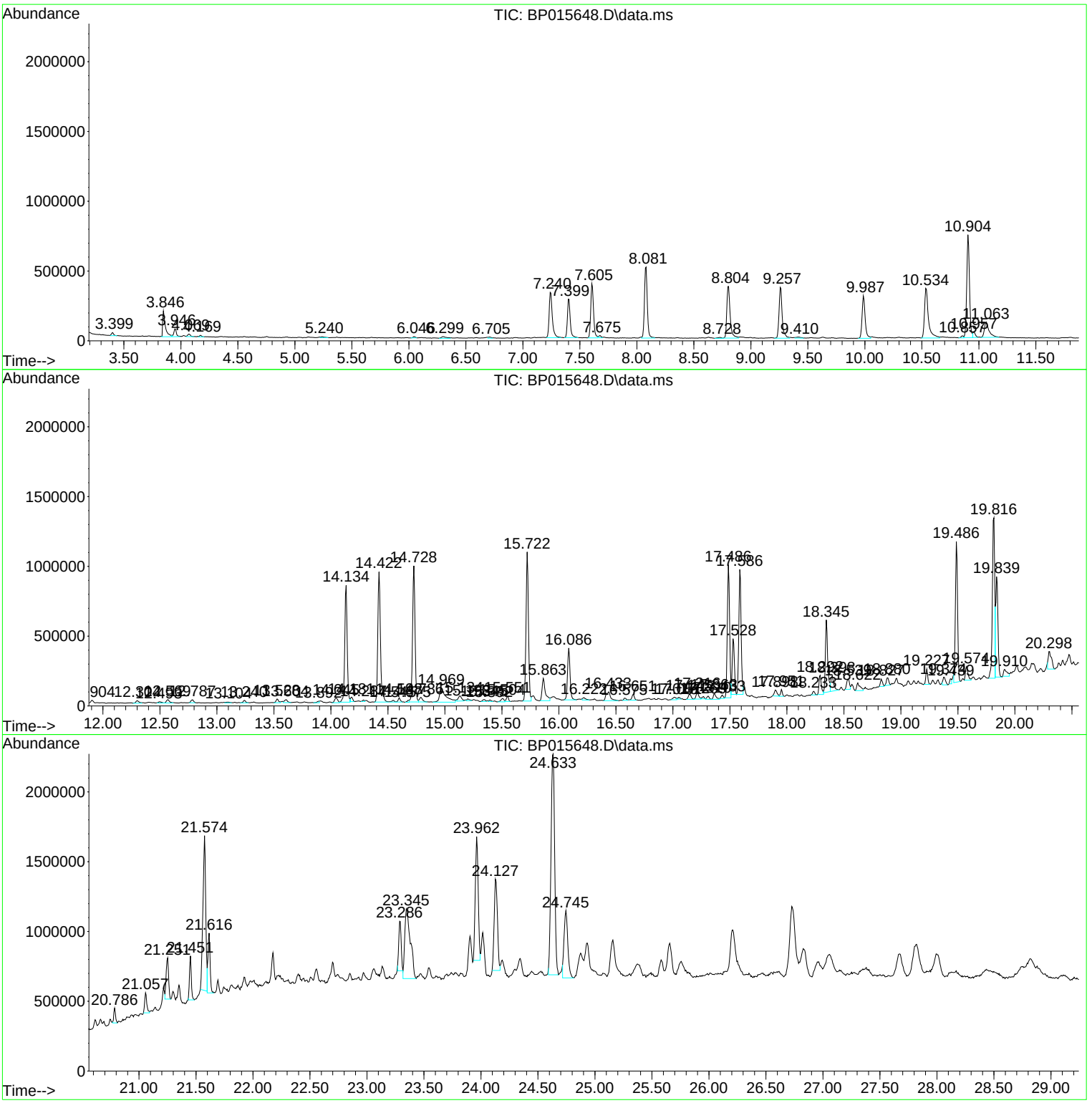
Instrument :
BNA_P
ClientSampleId :
DCFK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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\BP062223\
Data File : BP015648.D
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

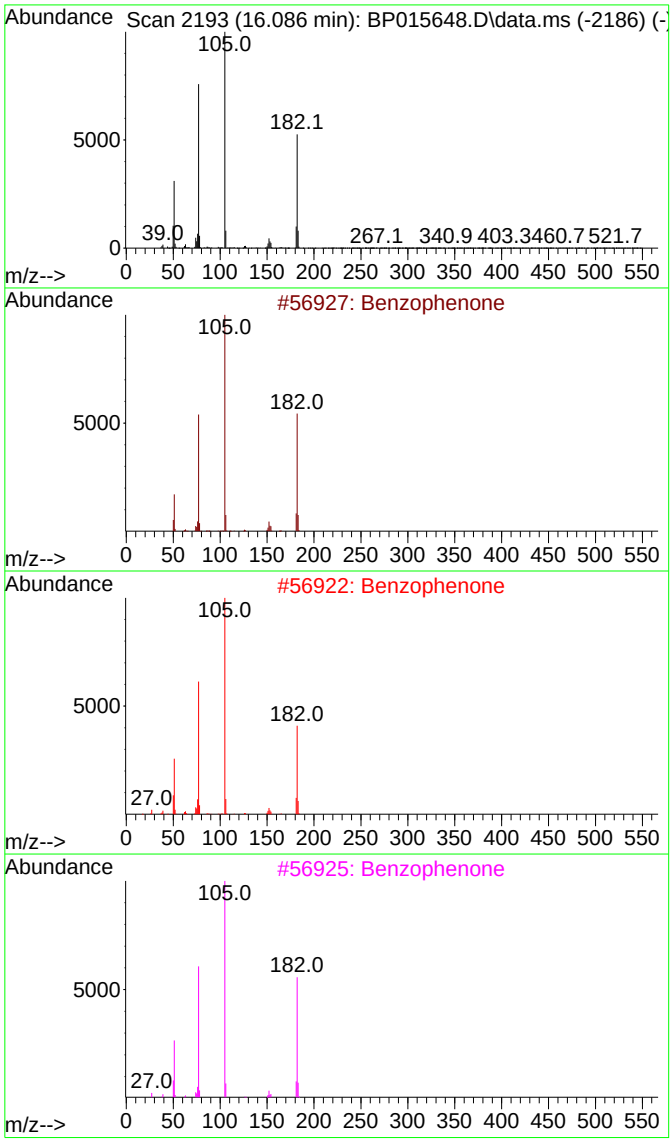
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

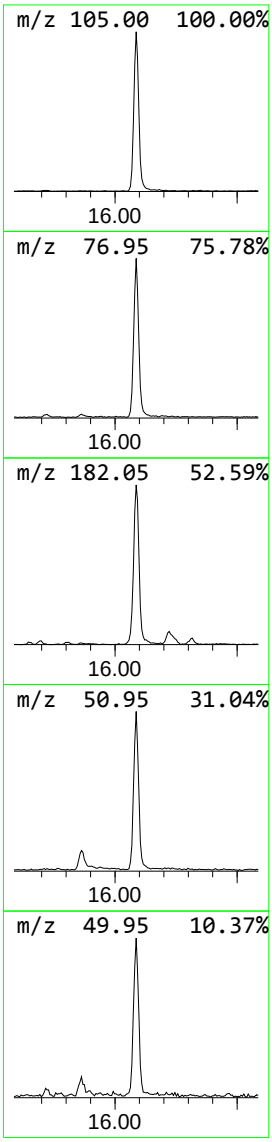
****

Peak Number 1 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.086	7.55 ng/ul	557616	Acenaphthene-d10	14.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	96
4	Benzophenone	182	C13H10O	000119-61-9	95
5	Benzophenone	182	C13H10O	000119-61-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015648.D
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Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

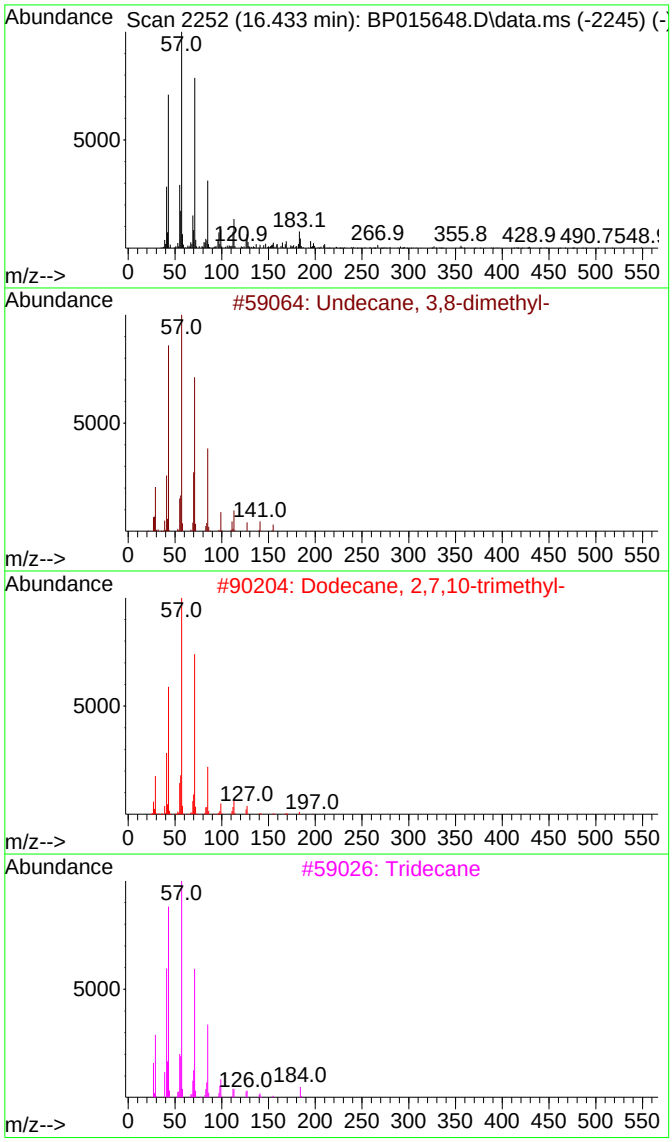
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

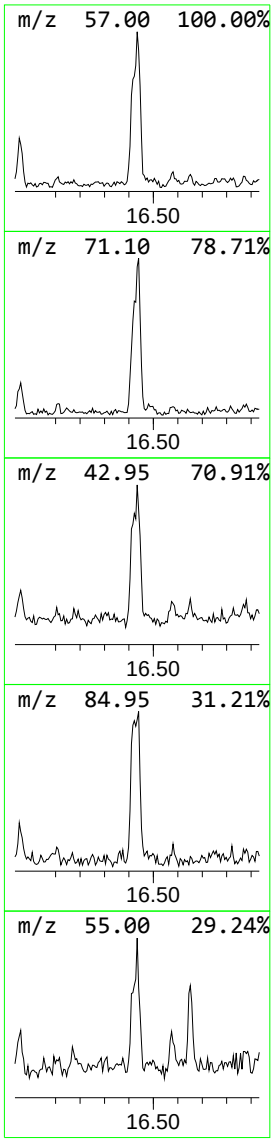
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.433	2.58 ng/ul	173574	Phenanthrene-d10	17.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	80
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
3	Tridecane	184	C13H28	000629-50-5	80
4	Tridecane	184	C13H28	000629-50-5	72
5	3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	72





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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

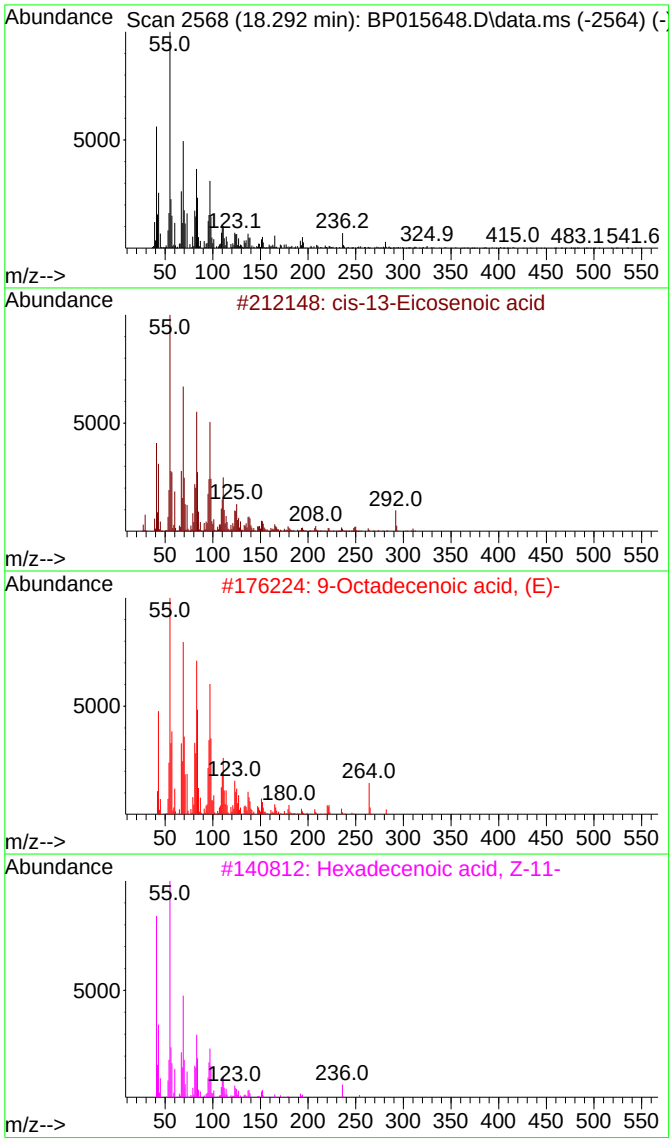
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

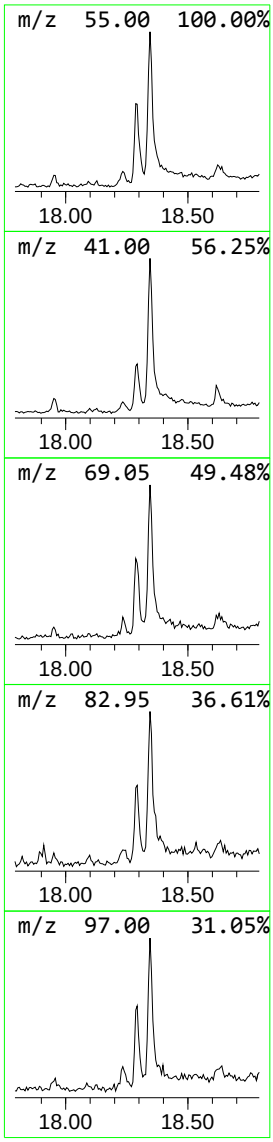
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 cis-13-Eicosenoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	3.03 ng/ul	203775	Phenanthrene-d10	17.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-13-Eicosenoic acid	310	C20H38O2	017735-94-3	99
2	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	94
3	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
4	6-Pentadecenoic acid, 13-methyl-...	254	C16H30O2	682751-34-4	91
5	Oleic Acid	282	C18H34O2	000112-80-1	91





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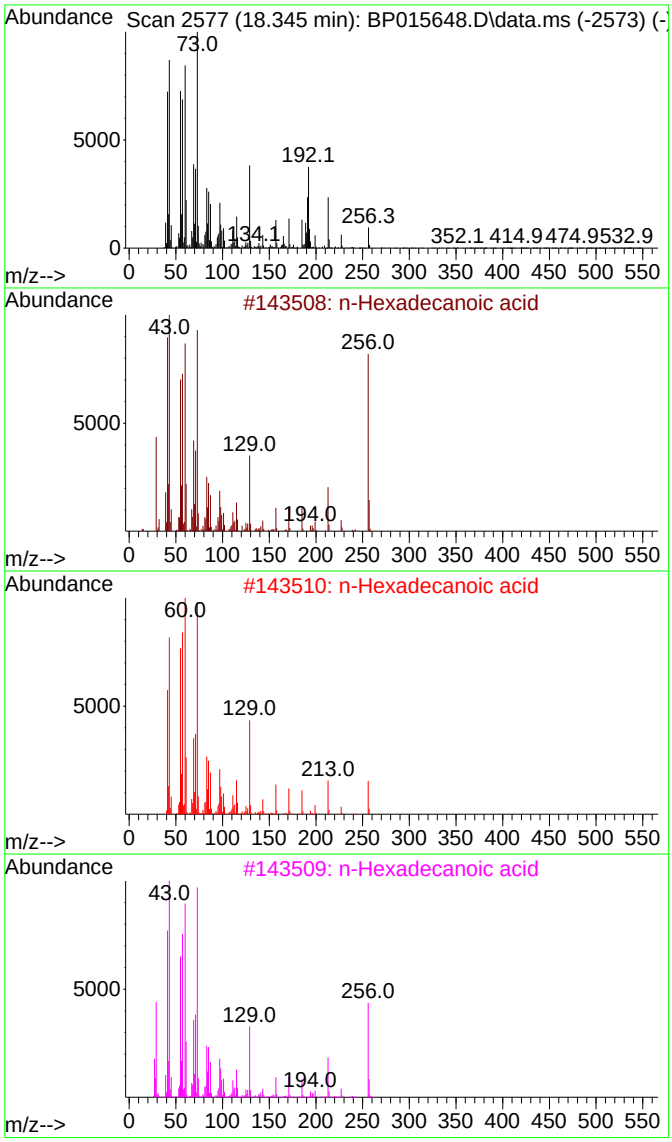
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

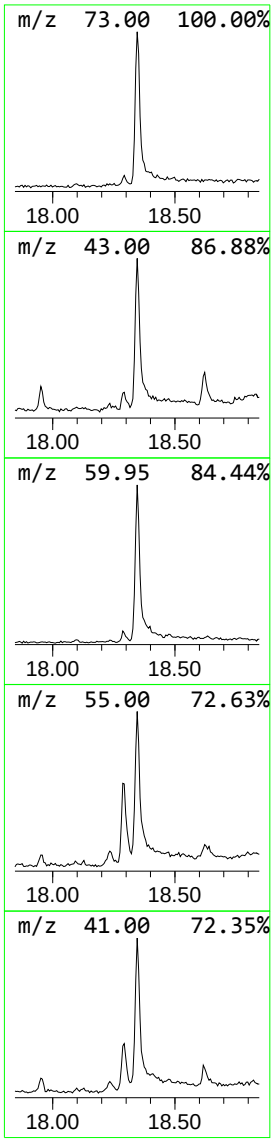
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	10.99 ng/ul	738558	Phenanthrene-d10	17.486

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	97
4	Octadecanoic acid	284	C18H36O2	000057-11-4	93
5	Tridecanoic acid	214	C13H26O2	000638-53-9	91





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Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

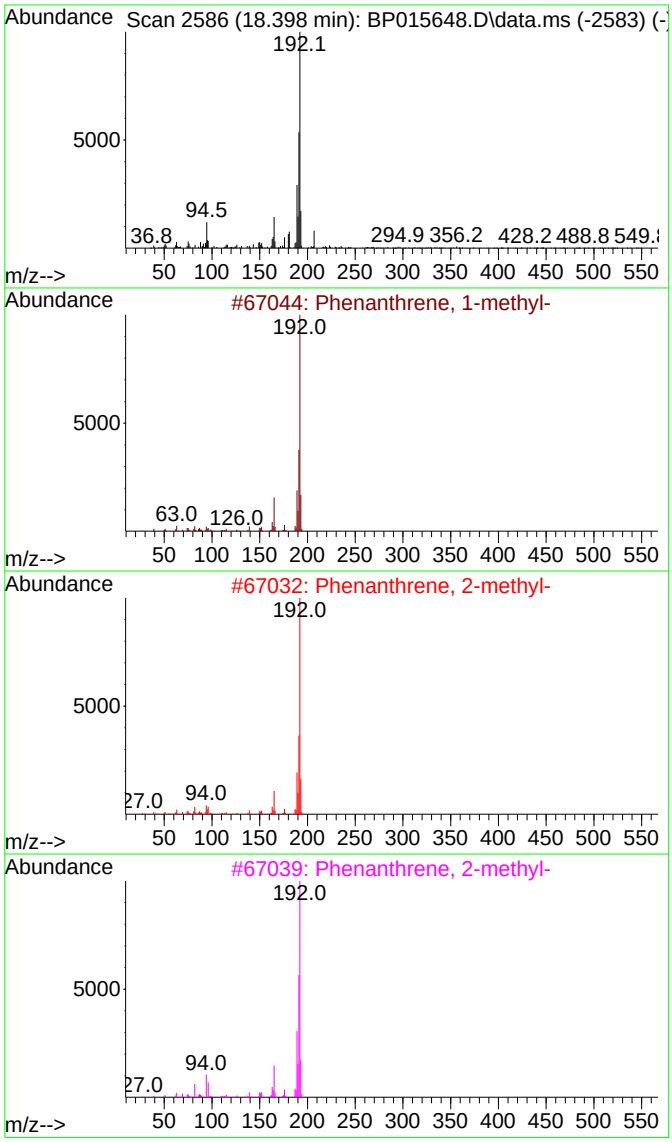
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

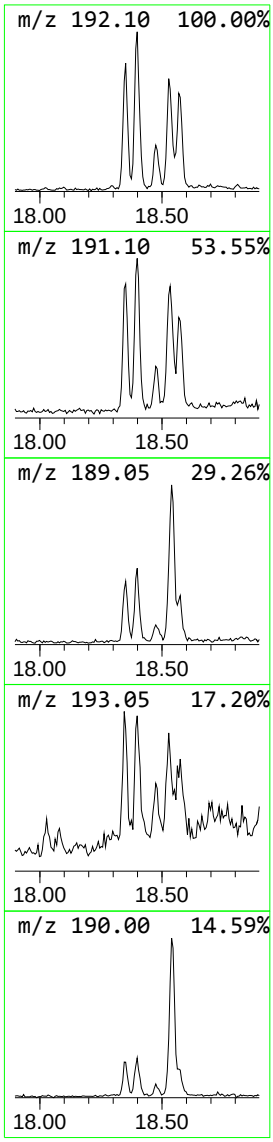
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	2.75 ng/ul	184636	Phenanthrene-d10	17.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90





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Data File : BP015648.D
Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

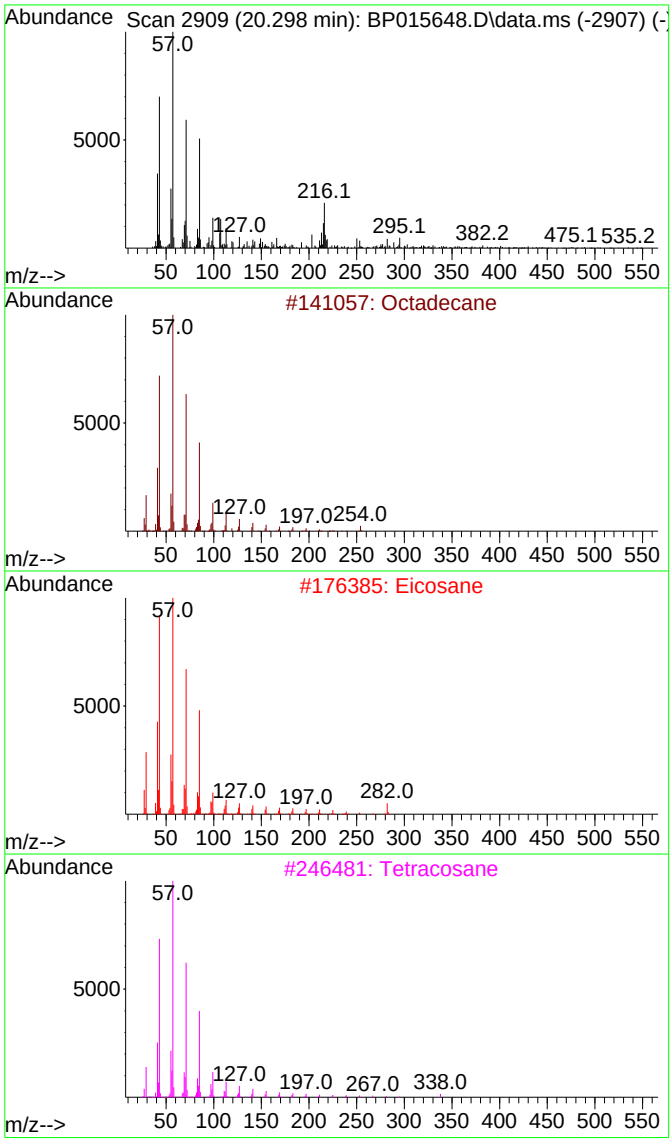
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

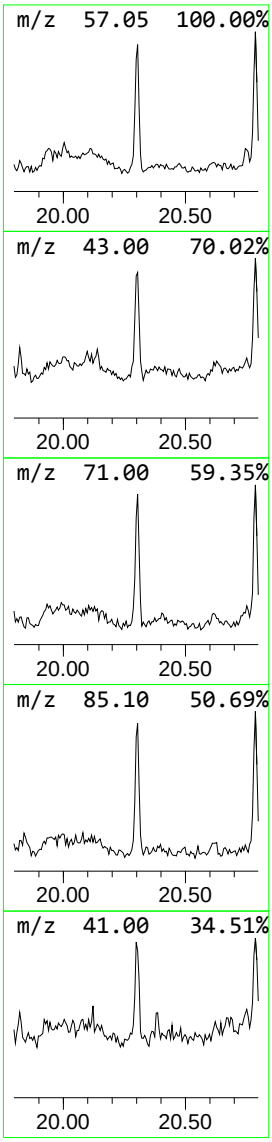
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.298	2.57 ng/ul	247306	Chrysene-d12	21.574

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	83
2	Eicosane	282	C20H42	000112-95-8	64
3	Tetracosane	338	C24H50	000646-31-1	58
4	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	58
5	Decane, 1-iodo-	268	C10H21I	002050-77-3	55





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Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

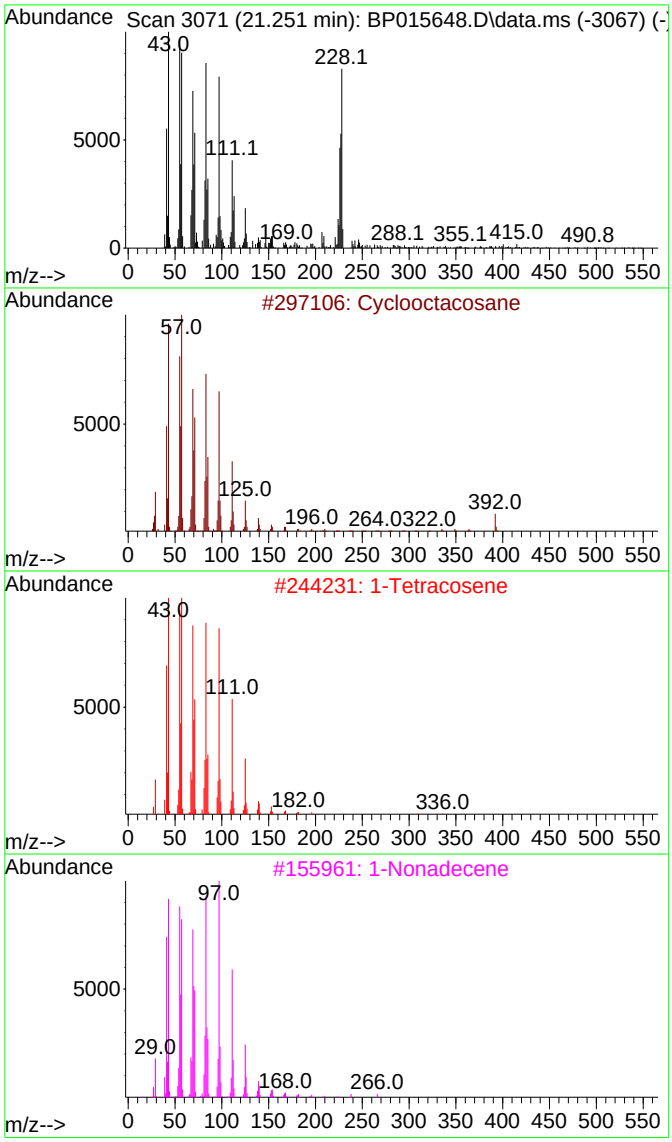
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

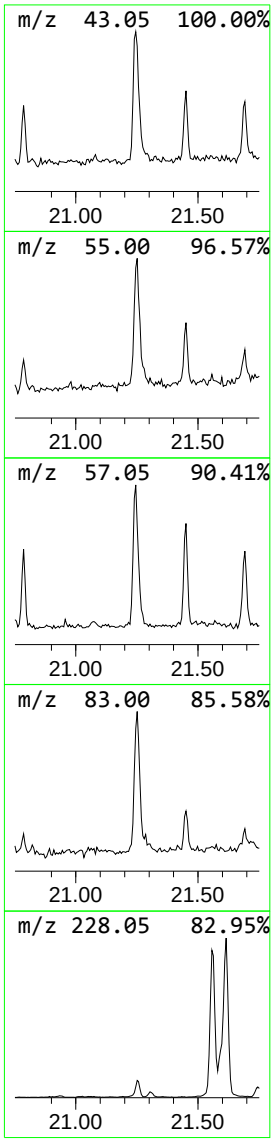
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic21.251 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	4.47 ng/ul	430184	Chrysene-d12	21.574

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctacosane	392	C28H56	000297-24-5	90
2	1-Tetracosene	336	C24H48	010192-32-2	89
3	1-Nonadecene	266	C19H38	018435-45-5	89
4	Cyclohexadecane	224	C16H32	000295-65-8	83
5	1-Docosene	308	C22H44	001599-67-3	83





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Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

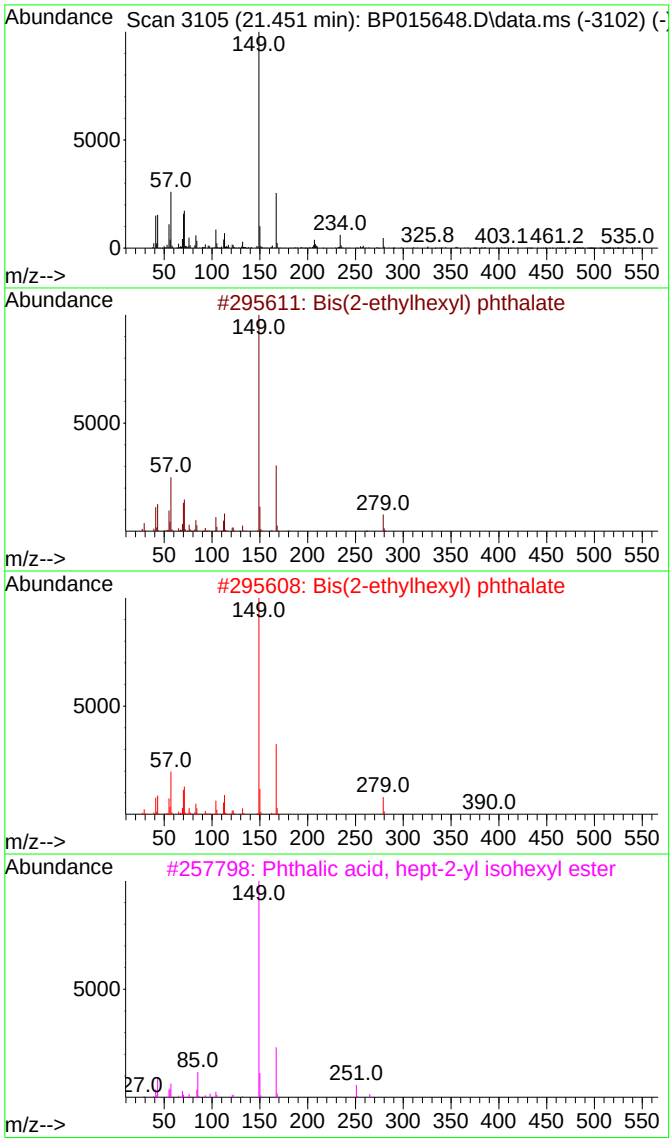
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

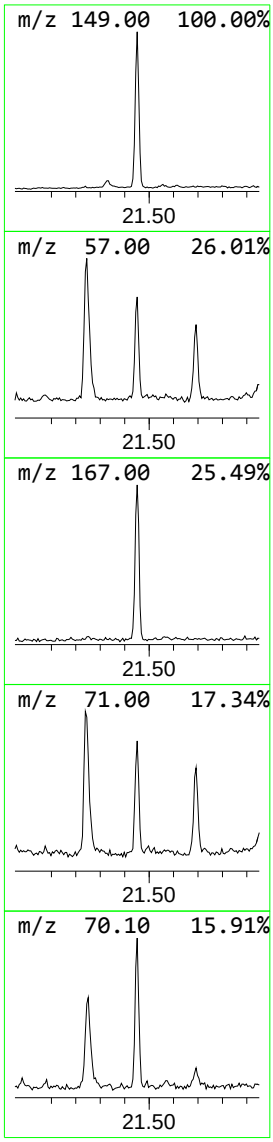
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Bis(2-ethylhexyl) phthalate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.451	3.61 ng/ul	347390	Chrysene-d12	21.574

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	87
2	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	74
3	Phthalic acid, hept-2-yl isohexy...	348	C21H32O4	1000356-97-3	64
4	Phthalic acid, dec-2-yl isohexyl...	390	C24H38O4	1000356-94-1	64
5	Phthalic acid, isohexyl 4-methyl...	334	C20H30O4	1010356-87-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

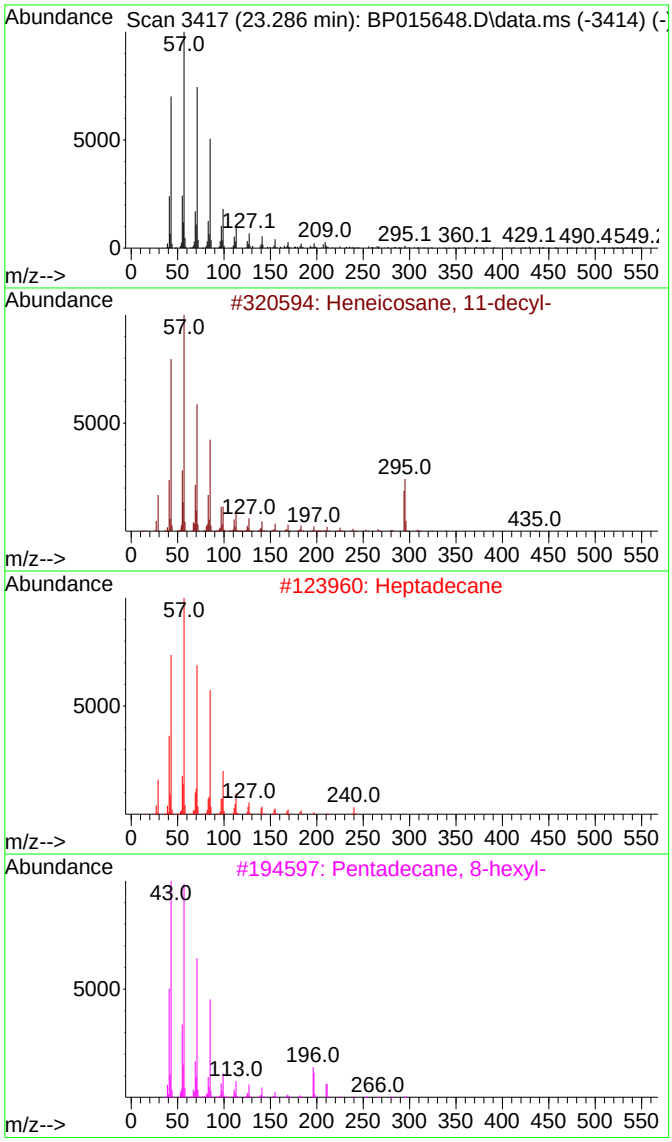
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

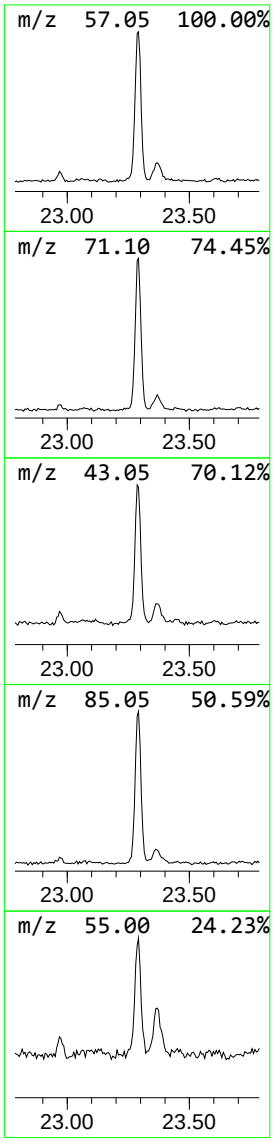
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.286	8.08 ng/ul	528405	Perylene-d12	24.127

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95
2	Heptadecane	240	C17H36	000629-78-7	94
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Docosane, 9-butyl-	366	C26H54	055282-14-9	91
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

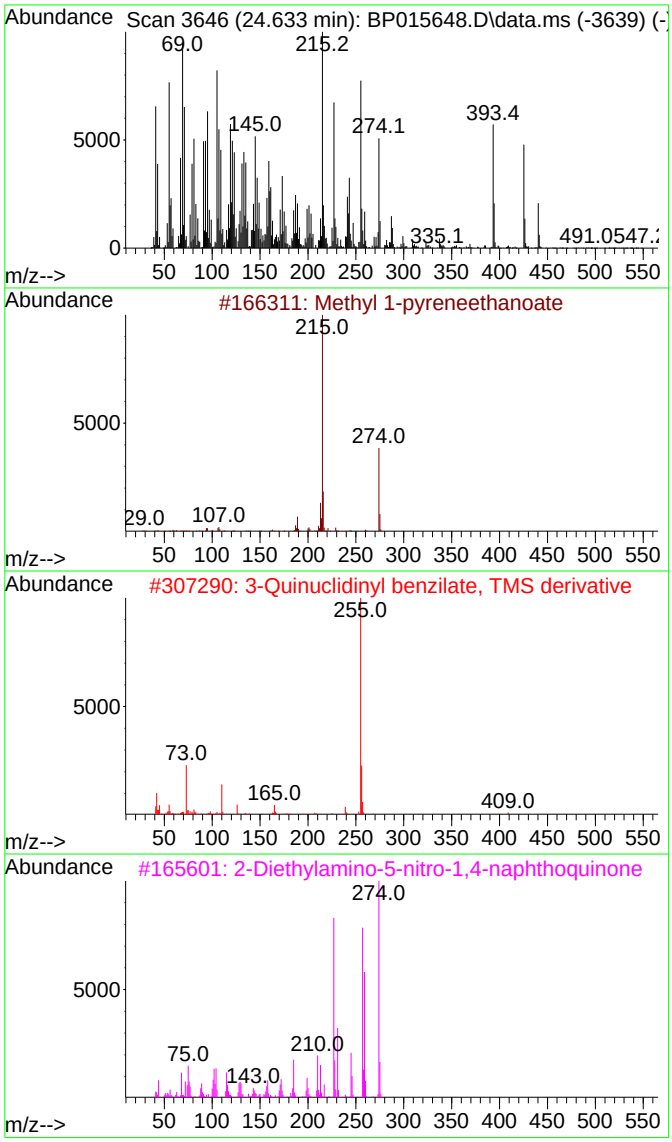
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

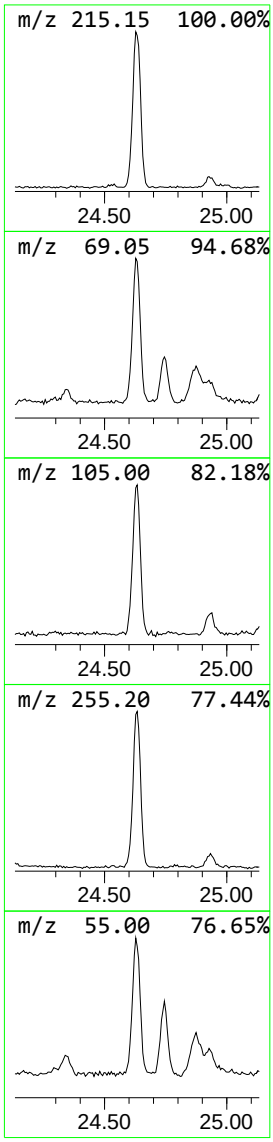
****

Peak Number 10 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	53.72 ng/ul	3515350	Perylene-d12	24.127

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	47
2	3-Quinuclidinyl benzilate, TMS d...	409	C24H31NO3Si	069569-91-1	25
3	2-Diethylamino-5-nitro-1,4-napht...	274	C14H14N2O4	110574-69-1	15
4	2-(4-Methyl-benzylsulfanyl)-3H-i...	255	C14H13N3S	1000296-58-2	15
5	Indanidine	215	C11H13N5	085392-79-6	15





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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

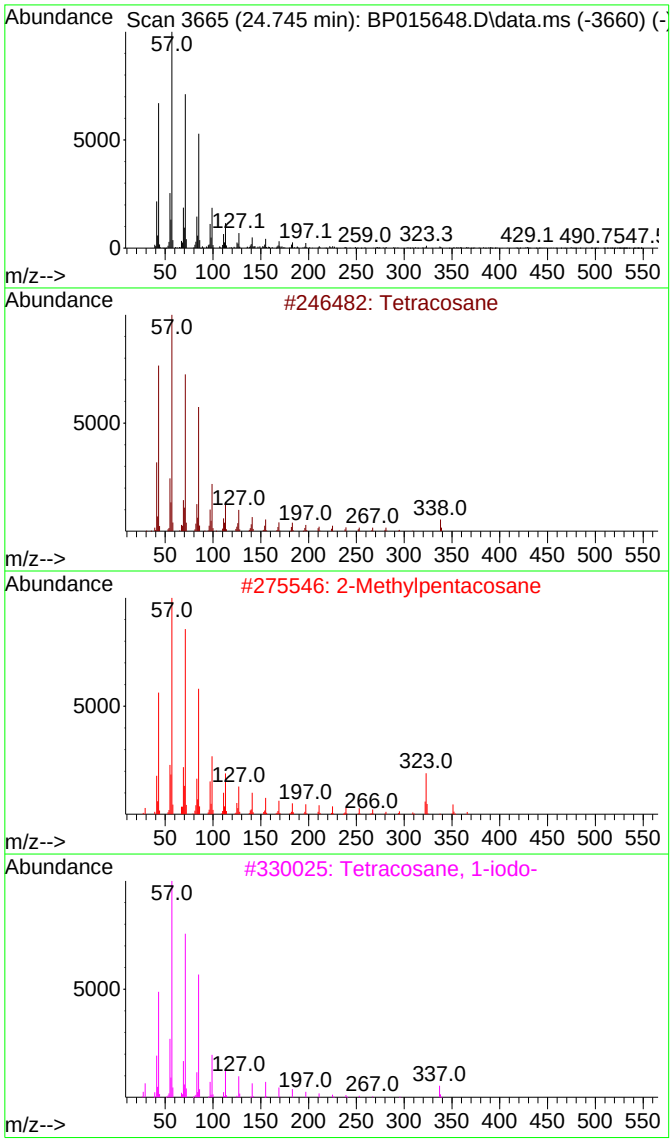
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

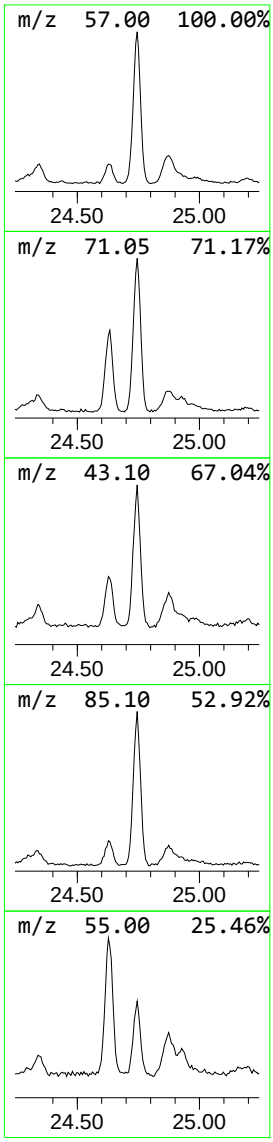
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.745	16.72 ng/ul	1094100	Perylene-d12	24.127

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	95
2	2-Methylpentacosane	366	C26H54	000629-87-8	93
3	Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	90
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015648.D
Acq On : 22 Jun 2023 15:39
Operator : MA/JU
Sample : 03083-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
Benzophenone	16.086	7.5	ng/ul	557616	3	14.728	1477660	20.0
(DEL) Alkane: S...	16.433	2.6	ng/ul	173574	4	17.486	1344380	20.0
cis-13-Eicoseno...	18.292	3.0	ng/ul	203775	4	17.486	1344380	20.0
n-Hexadecanoic ...	18.345	11.0	ng/ul	738558	4	17.486	1344380	20.0
Phenanthrene, 1...	18.398	2.8	ng/ul	184636	4	17.486	1344380	20.0
(DEL) Alkane: S...	20.298	2.6	ng/ul	247306	5	21.574	1926140	20.0
(DEL) Alkane: C...	21.251	4.5	ng/ul	430184	5	21.574	1926140	20.0
Bis(2-ethylhexy...	21.451	3.6	ng/ul	347390	5	21.574	1926140	20.0
(DEL) Alkane: S...	23.286	8.1	ng/ul	528405	6	24.127	1308730	20.0
unknown-01	24.633	53.7	ng/ul	3515350	6	24.127	1308730	20.0
(DEL) Alkane: S...	24.745	16.7	ng/ul	1094100	6	24.127	1308730	20.0