LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

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\SOM-EPA-BM061419MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	21	rVB	3817404	4965187	61.47%	5.111%
2	3.287	48	51	59	rVB	63999	86247	1.07%	0.089%
3	3.910	154	157	163	rVB	35652	49448	0.61%	0.051%
4	4.422	239	244	252	rVB	440617	578392	7.16%	0.595%
5	4.904	322	326	330	rBV	29568	53250	0.66%	0.055%
6	5.010	341	344	349	rVB	42051	51923	0.64%	0.053%
7	6.457	586	590	595	rVB4	11059	20232	0.25%	0.021%
8	6.769	639	643	647	rBV5	5937	10040	0.12%	0.010%
9	6.881	657	662	667	rBV	273652	379781	4.70%	0.391%
10	6.946	667	673	689	rVB	1332838	2221040	27.50%	2.286%
11	7.116	695	702	713	rBV	1094546	1695994	21.00%	1.746%
12	7.204	714	717	723	rVB3	13909	21783	0.27%	0.022%
13	7.310	728	735	745	rBV	1401816	2256519	27.93%	2.323%
14	7.516	766	770	775	rBV3	10415	16110	0.20%	0.017%
15	7.663	790	795	798	rVB3	7355	10891	0.13%	0.011%
16	7.775	807	814	821	rBV	995766	1546522	19.15%	1.592%
17	7.822	821	822	827	rVB4	8761	8285	0.10%	0.009%
18	8.028	853	857	865	rVB5	6009	12063	0.15%	0.012%
19	8.481	925	934	939	rBV	1634664	2692666	33.33%	2.772%
20	8.534	939	943	962	rVB	685649	1266503	15.68%	1.304%
21	8.940	1004	1012	1025	rBV	1380440	2283000	28.26%	2.350%
22	9.039	1025	1029	1035	rVB6	5986	9095	0.11%	0.009%
23	9.169	1046	1051	1056	rBV3	12784	21963	0.27%	0.023%
24	9.310	1067	1075	1081	rBV2	156956	231832	2.87%	0.239%
25	9.651	1126	1133	1143	rBV	1217979	1977270	24.48%	2.035%
26	10.181	1216	1223	1238	rBV	1883084	3192823	39.53%	3.287%
27	10.563	1280	1288	1296	rBV	1408933	2316858	28.68%	2.385%
28	10.704	1304	1312	1329	rBV	1489256	2676299	33.13%	2.755%
29	11.345	1418	1421	1428	rVB4	5769	10889	0.13%	0.011%
30	11.598	1457	1464	1469	rBV3	7573	16404	0.20%	0.017%
31	11.798	1489	1498	1504	rVB2	112217	198665	2.46%	0.205%
32	12.098	1538	1549	1554	rBV	56293	100460	1.24%	0.103%
33	12.157	1554	1559	1566	rVB	33503	57024	0.71%	0.059%
34	12.286	1571	1581	1585	rBV6	13396	25460	0.32%	0.026%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

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\SOM-EPA-BM061419MA.M
Title : SVOA CALIBRATION

35	12.375	1589	1596	1608	rBV	163957	315581	3.91%	0.325%
36	12.580	1627	1631	1634	rBV4	7241	9632	0.12%	0.010%
37	12.757	1658	1661	1667	rVB4	6973	10590	0.13%	0.011%
38	13.016	1701	1705	1709	rVB4	7258	10638	0.13%	0.011%
39	13.563	1793	1798	1808	rBV	42562	73834	0.91%	0.076%
40	13.827	1833	1843	1848	rBV	3685816	5159199	63.87%	5.311%
41	13.874	1848	1851	1859	rVB	1227665	1563014	19.35%	1.609%
42	14.110	1884	1891	1900	rVV	4180501	5946415	73.61%	6.121%
43	14.416	1936	1943	1951	rBV2	2244016	3307638	40.95%	3.405%
44	14.621	1970	1978	1994	rBV	1148160	1816378	22.49%	1.870%
45	14.845	2009	2016	2021	rVB4	19659	34044	0.42%	0.035%
46	15.210	2072	2078	2082	rBV	47404	63908	0.79%	0.066%
47	15.368	2100	2105	2106	rBV	50127	54704	0.68%	0.056%
48	15.410	2106	2112	2120	rVB	5136306	7409129	91.72%	7.627%
49	15.533	2127	2133	2148	rBV	1343563	1856919	22.99%	1.912%
50	15.651	2148	2153	2160	rVB	61727	88438	1.09%	0.091%
51	15.874	2185	2191	2202	rBV7	25109	49454	0.61%	0.051%
52	16.121	2229	2233	2237	rBV6	6870	11717	0.15%	0.012%
53	16.192	2241	2245	2251	rVB7	12952	17336	0.21%	0.018%
54	16.304	2259	2264	2269	rBV5	7545	10859	0.13%	0.011%
55	16.615	2313	2317	2324	rVB2	37019	48532	0.60%	0.050%
56	16.915	2363	2368	2370	rBV3	7301	9203	0.11%	0.009%
57	16.951	2370	2374	2379	rVB	10411	16010	0.20%	0.016%
58	17.162	2404	2410	2420	rBV2	2645642	3789568	46.91%	3.901%
59	17.262	2420	2427	2440	rVV	5471385	7509871	92.97%	7.731%
60	17.445	2456	2458	2464	rVB4	7821	12056	0.15%	0.012%
61	17.651	2488	2493	2495	rBV4	12129	15040	0.19%	0.015%
62	17.680	2495	2498	2504	rVB	22331	30962	0.38%	0.032%
63	18.051	2556	2561	2570	rBV	221643	322735	4.00%	0.332%
64	18.127	2570	2574	2583	rVB	210813	264590	3.28%	0.272%
65	18.345	2608	2611	2616	rVB3	17923	22127	0.27%	0.023%
66	18.627	2656	2659	2662	rVB3	21029	23191	0.29%	0.024%
67	18.974	2714	2718	2723	rBV	61839	84561	1.05%	0.087%
68	19.186	2750	2754	2758	rBV	71328	105651	1.31%	0.109%
69	19.327	2774	2778	2785	rBV2	58271	97031	1.20%	0.100%
70	19.439	2792	2797	2802	rVB2	48530	82476	1.02%	0.085%
71	19.556	2810	2817	2823	rBV	6180441	8077795	100.00%	8.316%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
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Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

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\SOM-EPA-BM061419MA.M
Title : SVOA CALIBRATION

72	20.092	2905	2908	2911	rBV	111728	109085	1.35%	0.112%
73	20.586	2988	2992	2996	rBV	158220	178423	2.21%	0.184%
74	21.050	3067	3071	3075	rBV	499128	574390	7.11%	0.591%
75	21.256	3103	3106	3111	rVB	76448	81465	1.01%	0.084%
76	21.345	3115	3121	3132	rVV	2893966	3572066	44.22%	3.677%
77	21.486	3142	3145	3151	rVB	295920	300213	3.72%	0.309%
78	21.927	3216	3220	3227	rBV	341917	436676	5.41%	0.450%
79	22.397	3296	3300	3305	rBV	310471	423444	5.24%	0.436%
80	22.909	3383	3387	3392	rBV	390759	504196	6.24%	0.519%
81	23.474	3476	3483	3494	rVB	4114236	7610116	94.21%	7.834%
82	23.615	3501	3507	3517	rVB2	1791111	3398120	42.07%	3.498%
83	24.139	3592	3596	3604	rVB	309914	570893	7.07%	0.588%

Sum of corrected areas: 97140831

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
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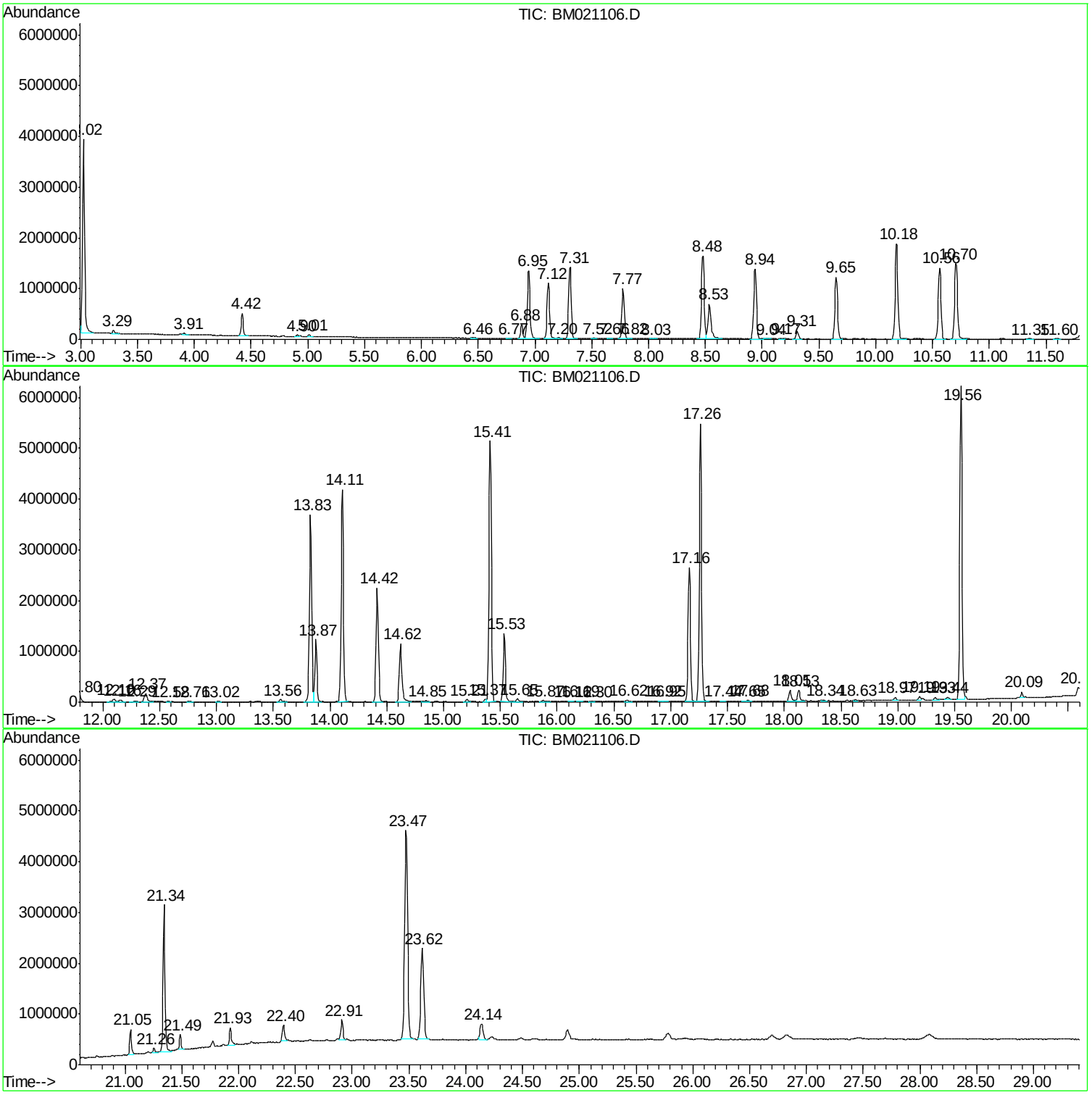
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

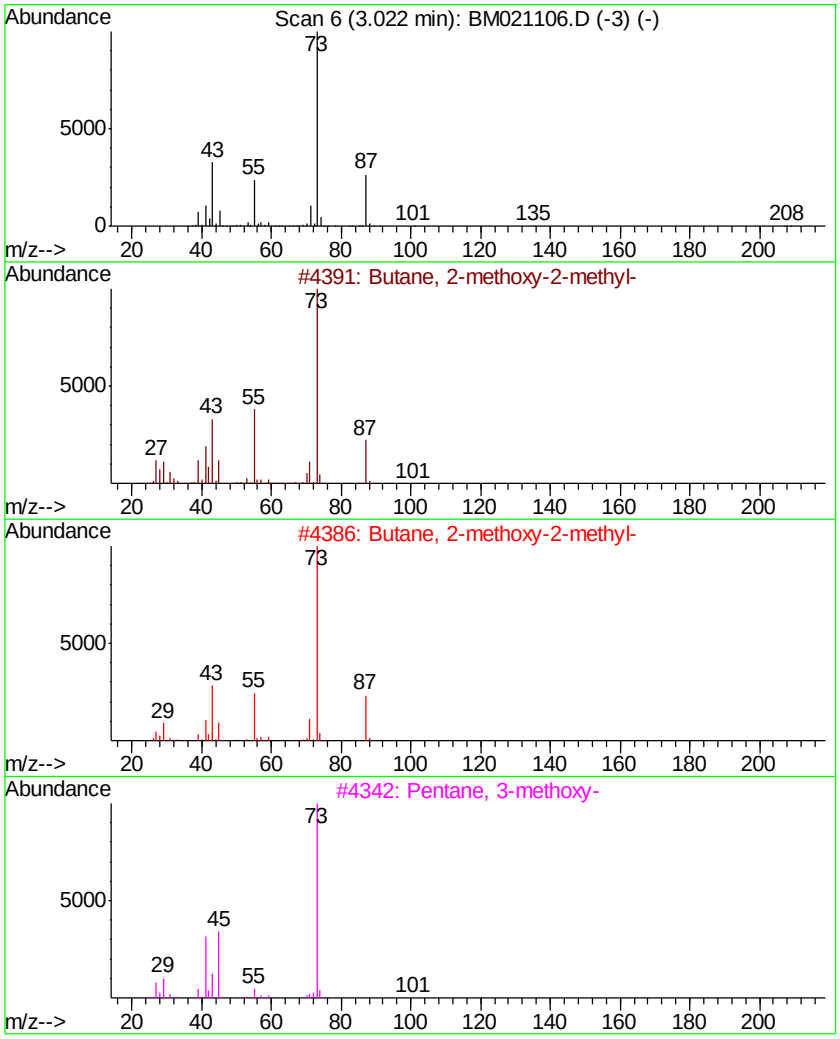
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

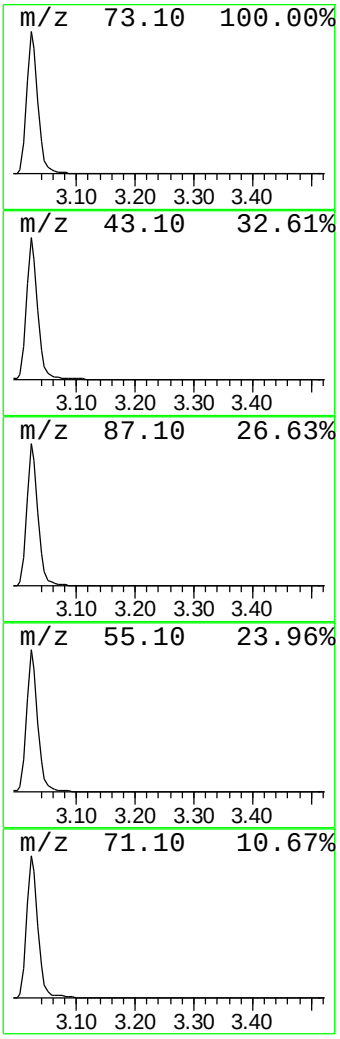
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	64.21 ng/ul	4965190	1,4-Dichlorobenzene-d4	7.77

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	78
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

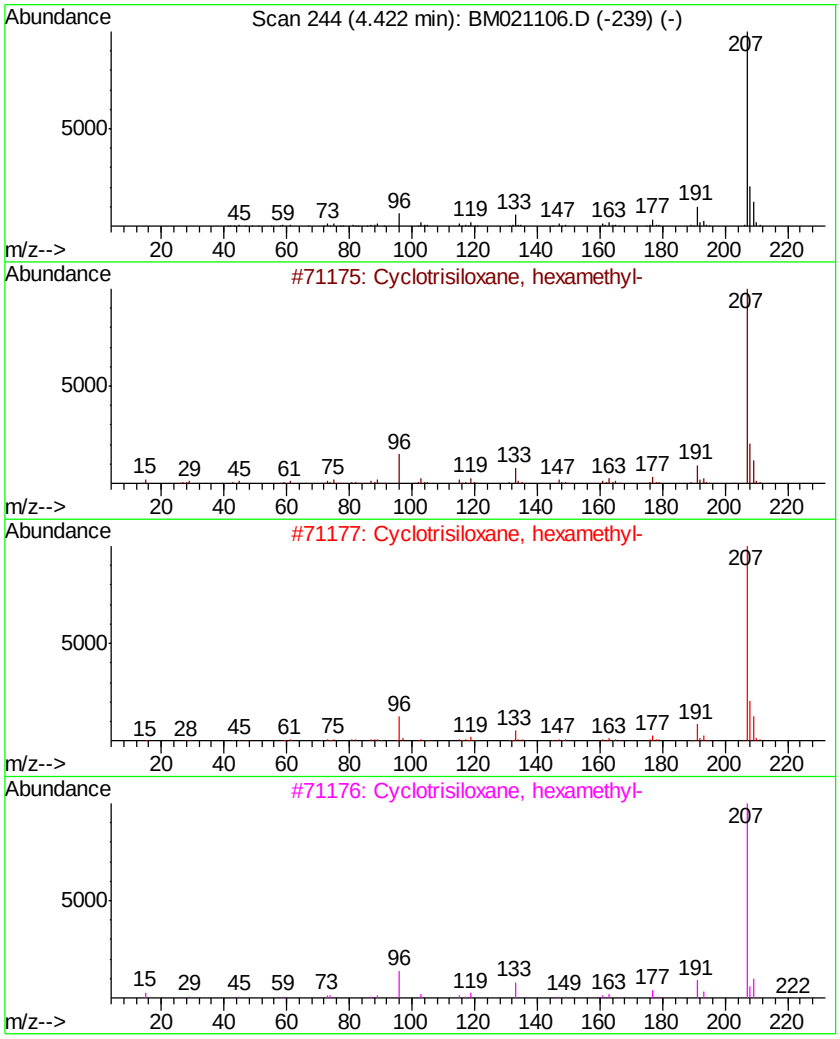
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

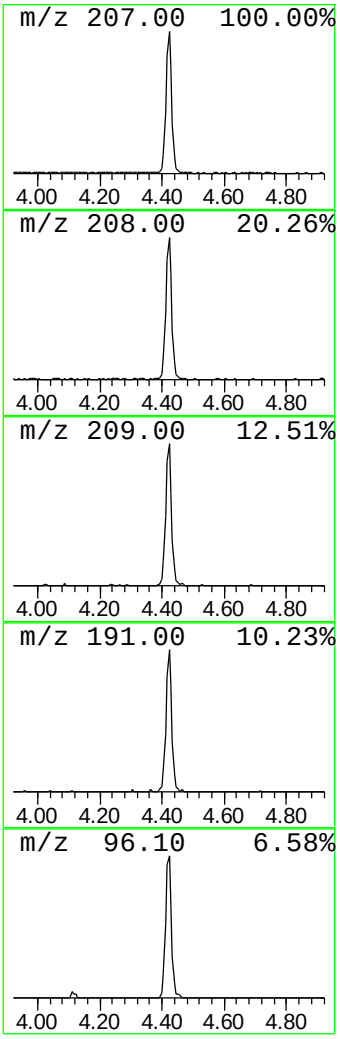
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclotrisiloxane, hexamethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.42	7.48 ng/ul	578392	1,4-Dichlorobenzene-d4	7.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	90
2	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	87
3	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	43
4	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	39





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

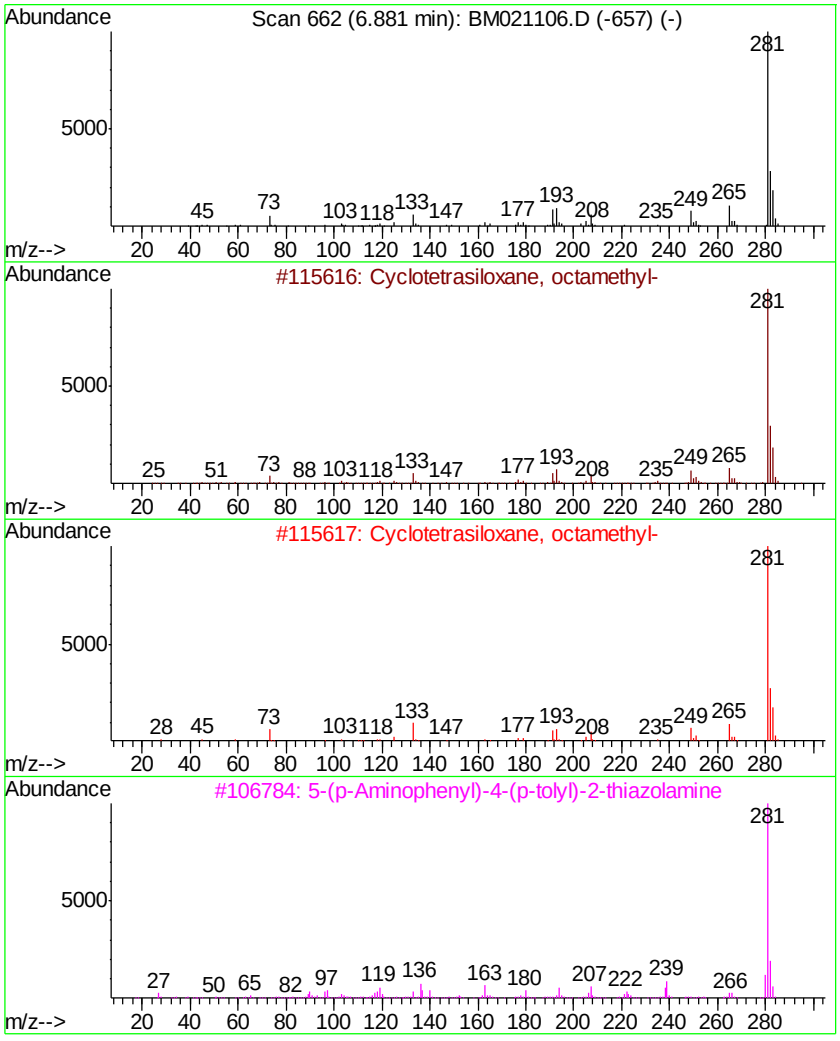
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

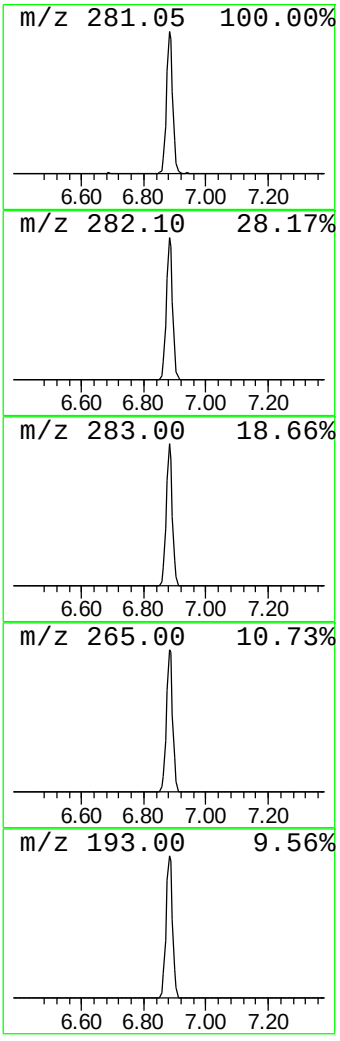
****

Peak Number 3 Cyclotetrasiloxane, octamet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	4.91 ng/ul	379781	1,4-Dichlorobenzene-d4	7.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	91
2	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	91
3	5-(p-Aminophenyl)-4-(p-tolyl)-2-...	281	C16H15N3S	094460-46-5	59
4	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	56
5	7H-Dibenzo[b,g]carbazole, 7-methyl-	281	C21H15N	003557-49-1	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

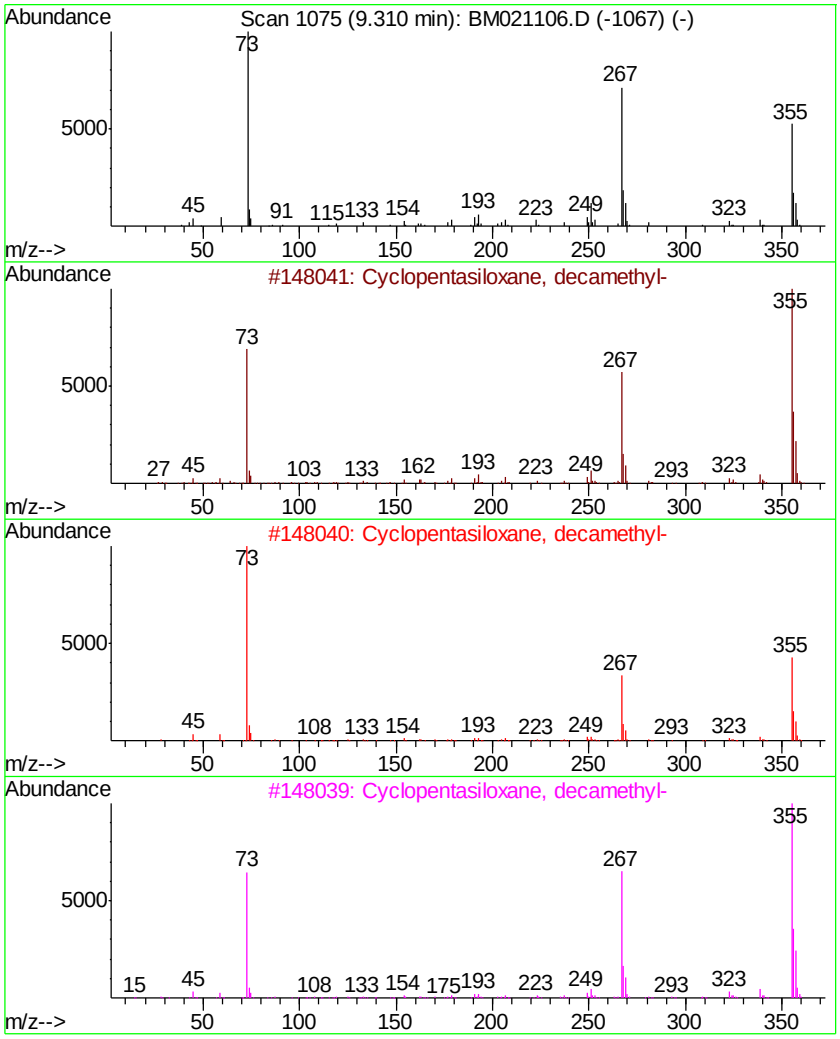
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

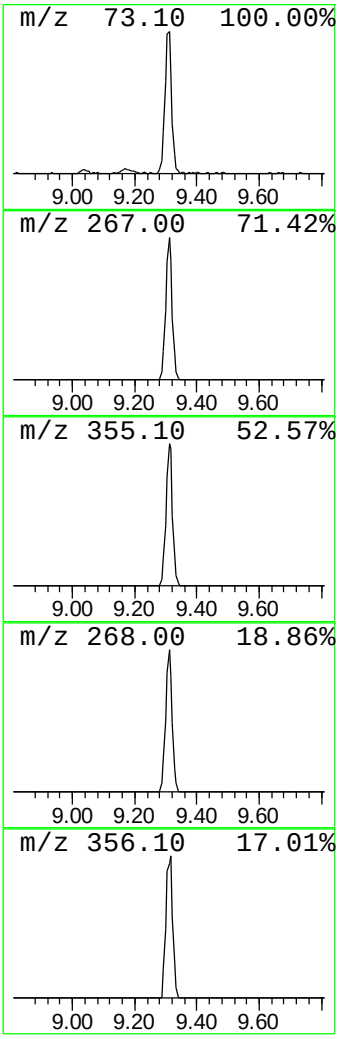
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentasiloxane, decamet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.31	2.00 ng/ul	231832	Naphthalene-d8	10.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	87
3	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	86
4	Isoproterenol tri-TMS derivative	427	C20H41NO3Si3	029522-12-1	43
5	3,4-Dihydroxymandelic acid, ethy...	428	C19H36O5Si3	1000071-70-2	43





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

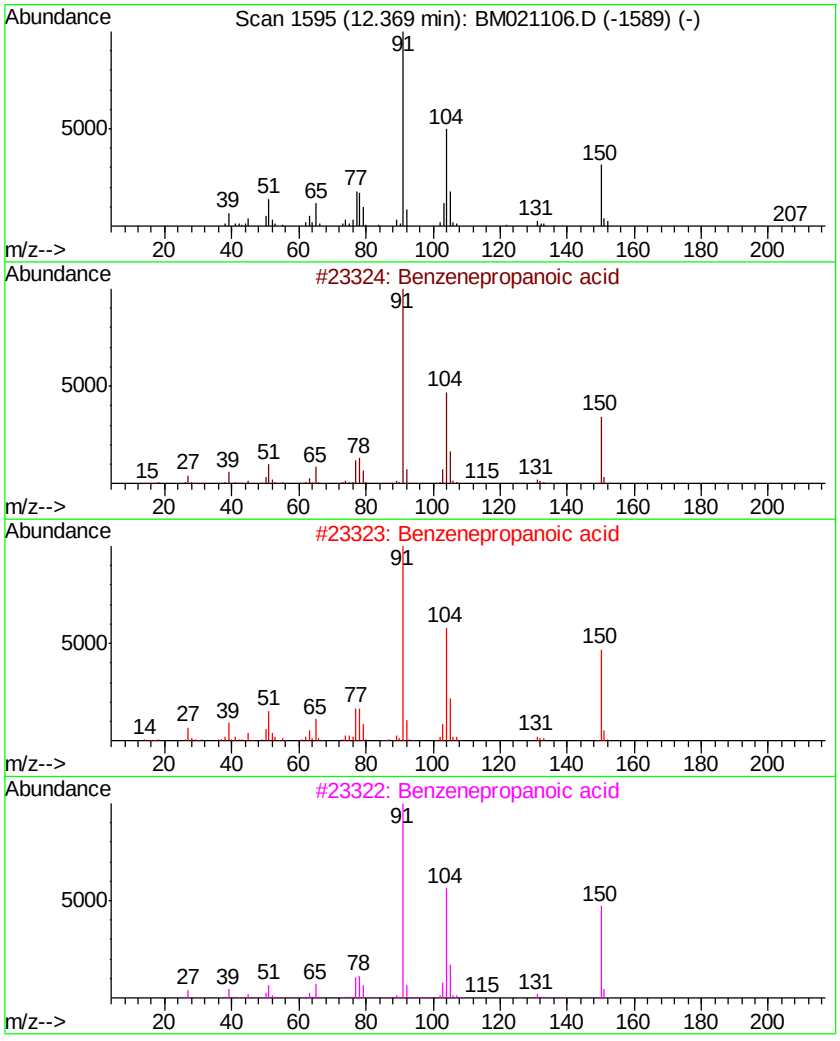
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

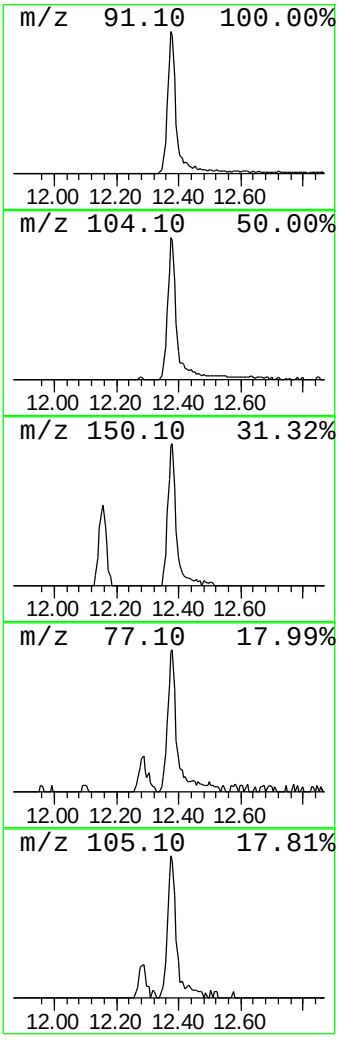
****

Peak Number 5 Benzenepropanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	2.72 ng/ul	315581	Naphthalene-d8	10.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenepropanoic acid	150	C9H10O2	000501-52-0	95
2	Benzenepropanoic acid	150	C9H10O2	000501-52-0	94
3	Benzenepropanoic acid	150	C9H10O2	000501-52-0	91
4	Benzylmalonic acid	194	C10H10O4	000616-75-1	91
5	4-Nitrosophenyl-.beta.-phenylpro...	255	C15H13NO3	1000129-40-0	78





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

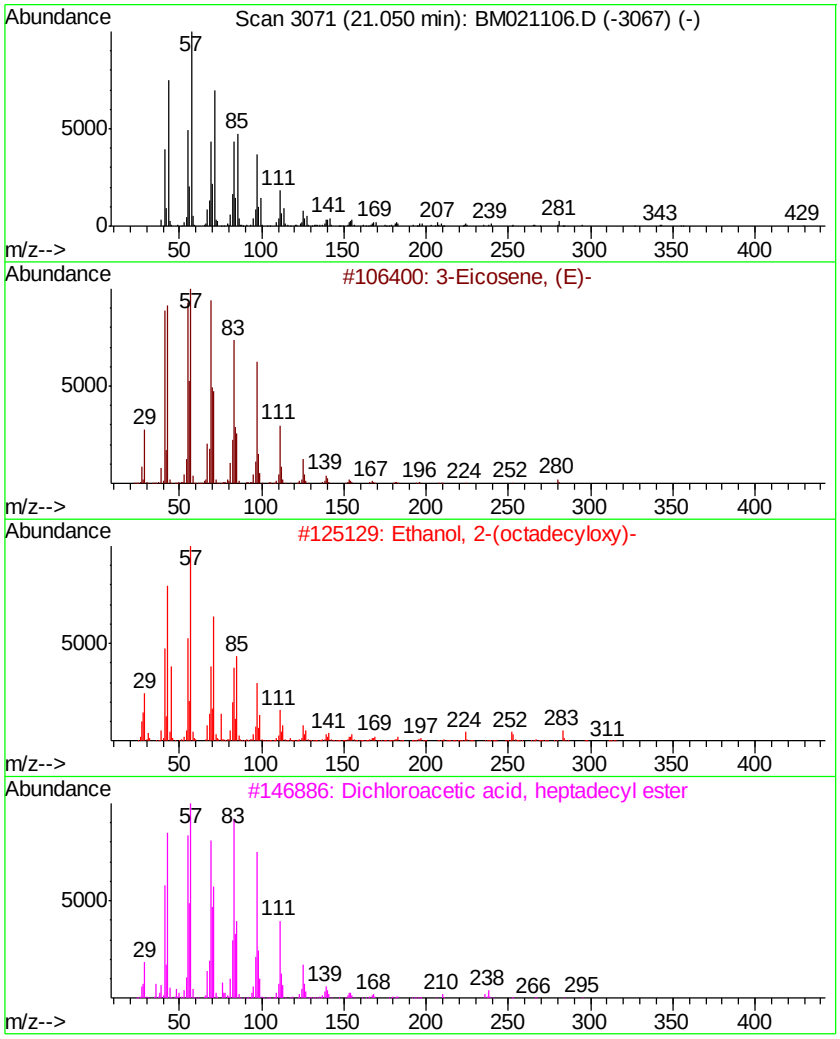
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

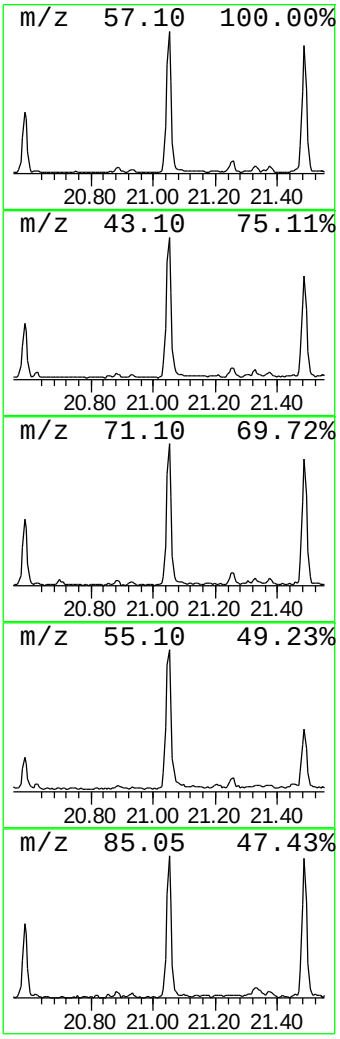
****

Peak Number 6 (DEL) Alkane: Cyclic21.05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.05	3.22 ng/ul	574390	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	90
2	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	87
3	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	70
4	1-Docosanethiol	342	C22H46S	007773-83-3	70
5	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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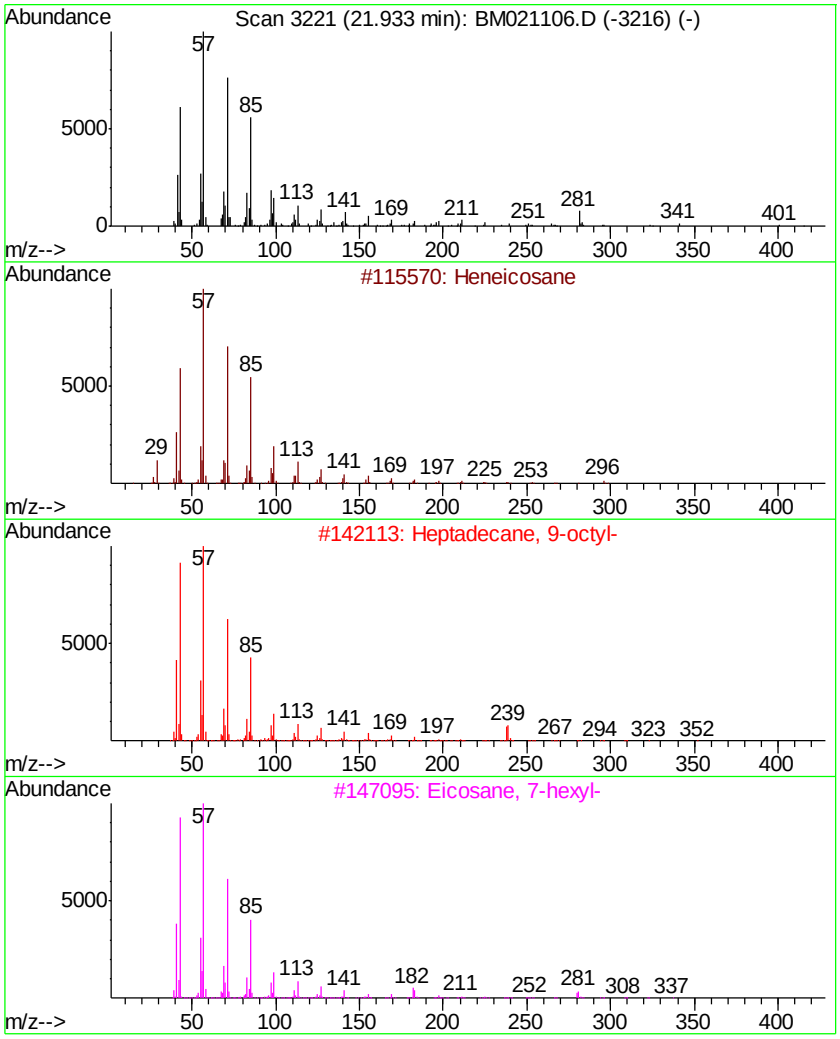
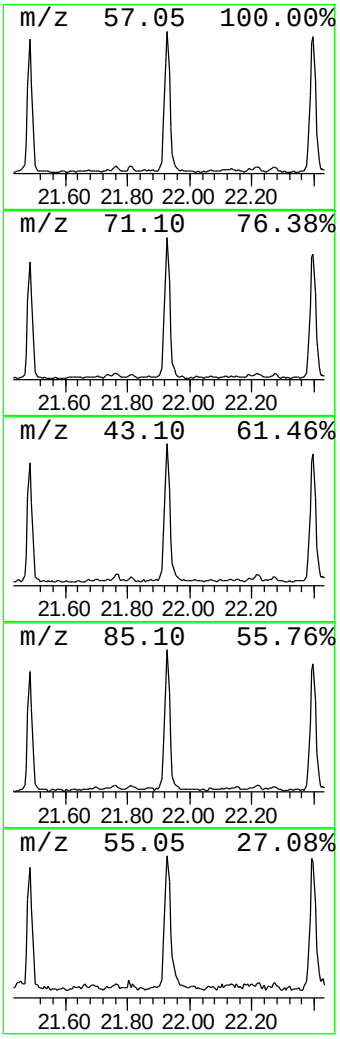
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.93	2.44 ng/ul	436676	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4	Docosane, 7-hexyl-	394	C28H58	055373-86-9	91
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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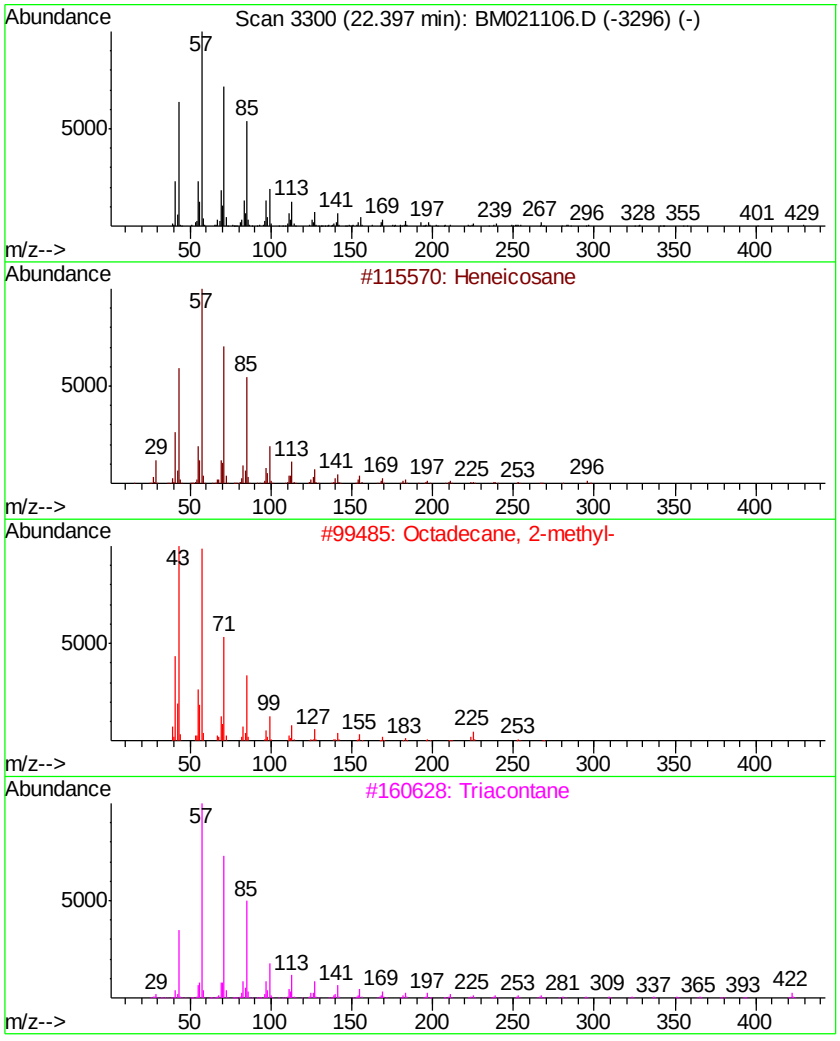
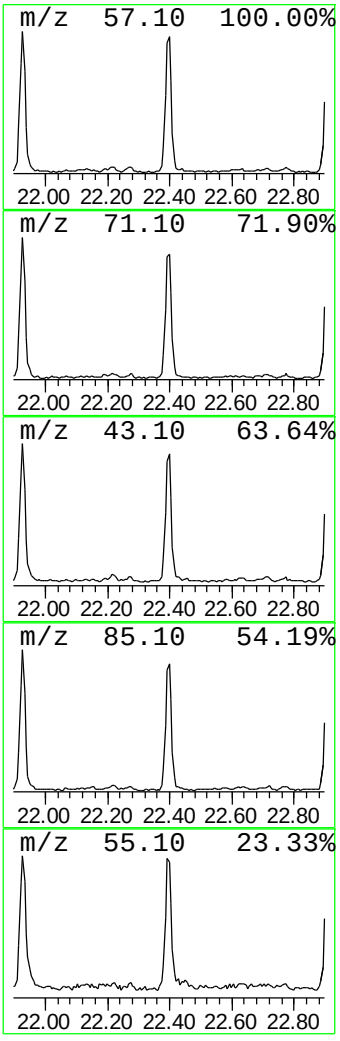
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.40	2.37 ng/ul	423444	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5	Tetracosane	338	C24H50	000646-31-1	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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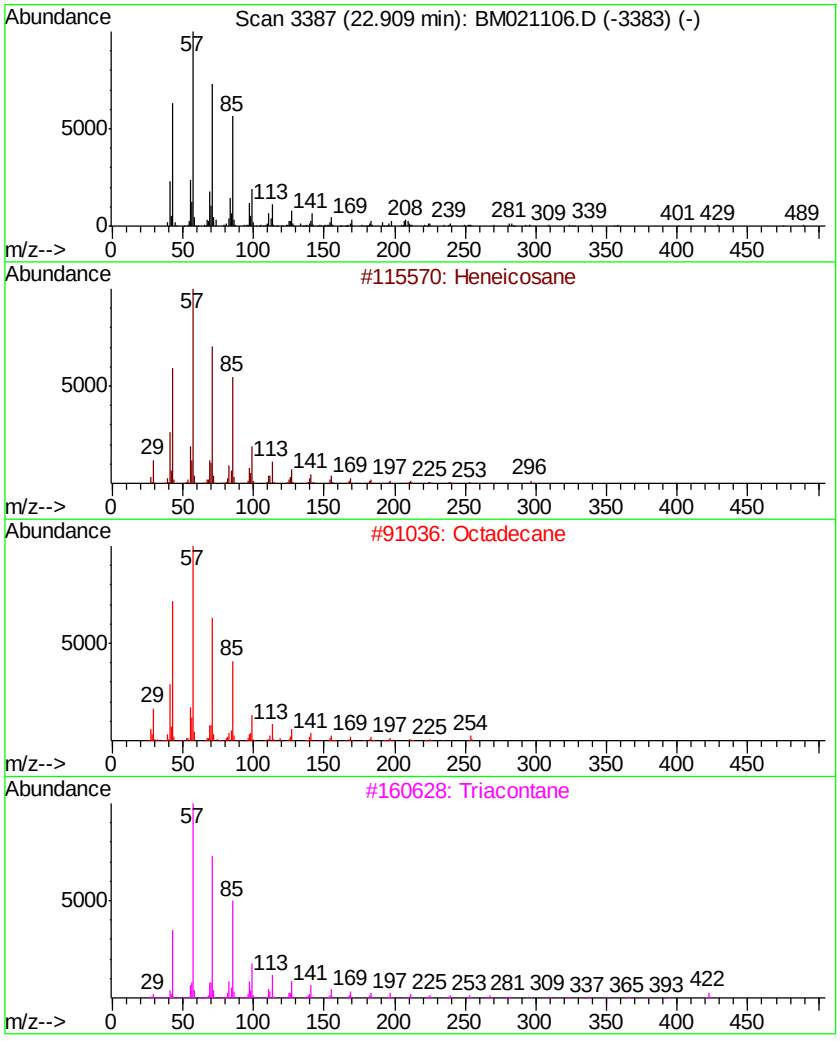
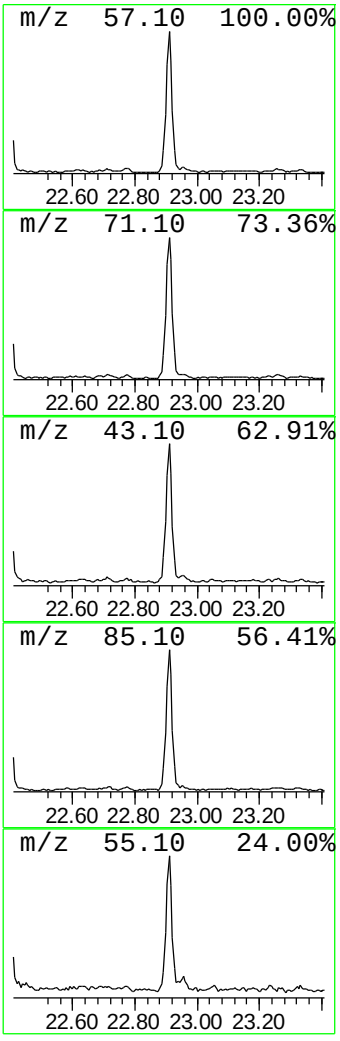
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.91	2.97 ng/ul	504196	Perylene-d12	23.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Octadecane	254	C18H38	000593-45-3	96
3	Triacontane	422	C30H62	000638-68-6	91
4	Nonacosane	408	C29H60	000630-03-5	91
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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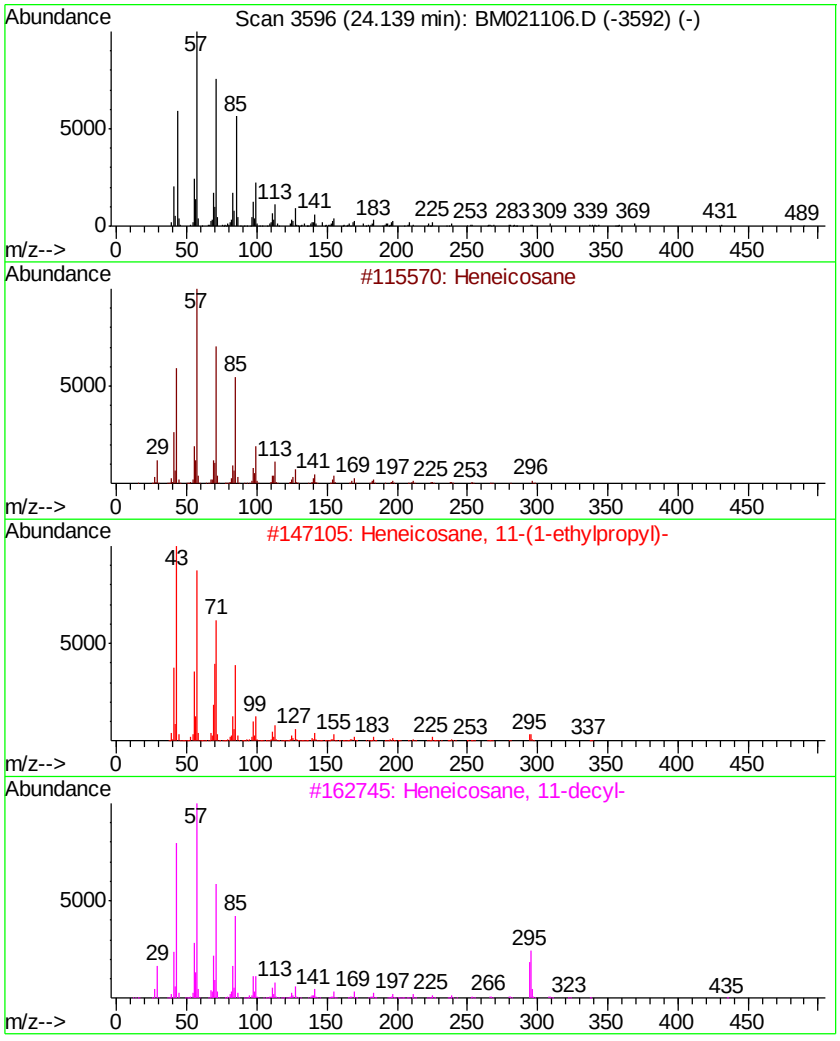
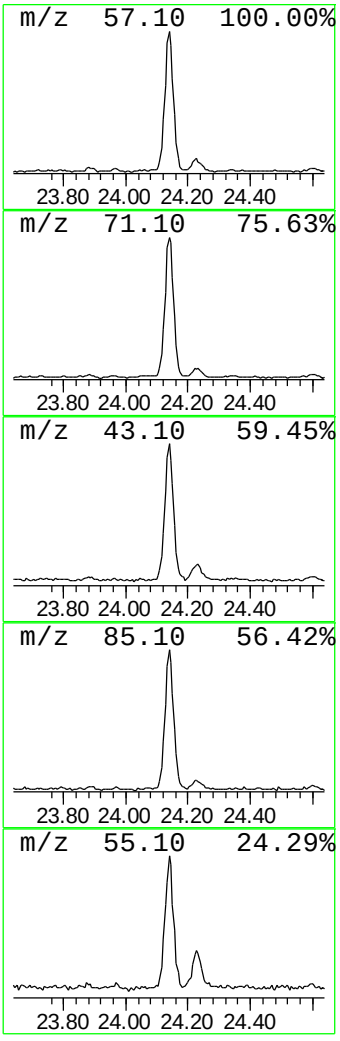
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.14	3.36 ng/ul	570893	Perylene-d12	23.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
4	Tetratetracontane	619	C44H90	007098-22-8	90
5	Nonacosane	408	C29H60	000630-03-5	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021106.D
Acq On : 22 Jun 2019 18:05
Operator : HP/JU
Sample : K3362-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A4200

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	64.2	ng/ul	4965190	1	7.77	1546520	20.0
Cyclotrisiloxane,...	4.42	7.5	ng/ul	578392	1	7.77	1546520	20.0
Cyclotetrasiloxan...	6.88	4.9	ng/ul	379781	1	7.77	1546520	20.0
Cyclopentasiloxan...	9.31	2.0	ng/ul	231832	2	10.56	2316860	20.0
Benzenepropanoic ...	12.37	2.7	ng/ul	315581	2	10.56	2316860	20.0
(DEL) Alkane: Cyc...	21.05	3.2	ng/ul	574390	5	21.34	3572070	20.0
(DEL) Alkane: Str...	21.93	2.4	ng/ul	436676	5	21.34	3572070	20.0
(DEL) Alkane: Str...	22.40	2.4	ng/ul	423444	5	21.34	3572070	20.0
(DEL) Alkane: Str...	22.91	3.0	ng/ul	504196	6	23.61	3398120	20.0
(DEL) Alkane: Str...	24.14	3.4	ng/ul	570893	6	23.61	3398120	20.0