LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

Signal : TIC: BN015105.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.281	30	33	41	rVB	31157	42242	0.49%	0.093%
2	3.546	76	78	83	rBV	6654	9784	0.11%	0.021%
3	3.676	97	100	112	rBV	270321	405985	4.74%	0.890%
4	3.781	117	118	123	rVB4	11161	10888	0.13%	0.024%
5	3.999	150	155	161	rVB	47319	68346	0.80%	0.150%
6	4.352	210	215	220	rBV2	18272	28377	0.33%	0.062%
7	4.487	233	238	242	rBV	13607	18674	0.22%	0.041%
8	4.534	242	246	252	rVV	45902	60038	0.70%	0.132%
9	5.281	369	373	380	rVB	35441	47402	0.55%	0.104%
10	5.411	389	395	405	rBV	102438	210349	2.46%	0.461%
11	5.770	451	456	462	rBV	51015	75795	0.89%	0.166%
12	5.846	464	469	475	rVB3	6286	10238	0.12%	0.022%
13	6.893	639	647	657	rBV	445308	696905	8.14%	1.527%
14	7.064	667	676	683	rBV	347316	544870	6.36%	1.194%
15	7.258	702	709	718	rBV	457086	721044	8.42%	1.580%
16	7.328	718	721	725	rBV	10878	16667	0.19%	0.037%
17	7.722	782	788	795	rBV	565314	907541	10.60%	1.989%
18	7.787	795	799	805	rVB2	12104	16823	0.20%	0.037%
19	7.964	825	829	837	rVB4	7356	12359	0.14%	0.027%
20	8.275	876	882	885	rBV5	5165	9197	0.11%	0.020%
21	8.416	899	906	913	rBV	520403	813967	9.50%	1.784%
22	8.534	922	926	932	rVB	18339	28958	0.34%	0.063%
23	8.816	967	974	976	rBV	11111	17160	0.20%	0.038%
24	8.869	976	983	992	rVV	381070	630704	7.36%	1.382%
25	8.975	996	1001	1007	rVV5	4647	8684	0.10%	0.019%
26	9.311	1054	1058	1068	rVB6	3983	8706	0.10%	0.019%
27	9.587	1098	1105	1113	rBV	267628	444259	5.19%	0.974%
28	10.116	1188	1195	1209	rVB	506802	814209	9.51%	1.784%
29	10.275	1216	1222	1239	rBV	989520	1526745	17.83%	3.346%
30	10.499	1253	1260	1268	rVB	775864	1278077	14.92%	2.801%
31	10.628	1275	1282	1290	rBV	476949	788900	9.21%	1.729%
32	12.087	1524	1530	1536	rBV2	7461	12236	0.14%	0.027%
33	12.710	1631	1636	1642	rBV3	17402	25029	0.29%	0.055%
34	13.187	1711	1717	1723	rBV2	15385	21563	0.25%	0.047%
35	13.257	1723	1729	1736	rVV	201968	315526	3.68%	0.691%
36	13.557	1775	1780	1785	rBV	12259	16540	0.19%	0.036%

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Instrument : BNA_N
ClientSampleId : BGDB6

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_N\METHODS\SFAM-EPA-BN061821MA.M
Title : SVOA CALIBRATION

37	13.757	1808	1814	1821	rBV	835189	1147080	13.39%	2.514%
38	13.893	1833	1837	1842	rVB6	4894	8932	0.10%	0.020%
39	14.040	1851	1862	1869	rBV	1020503	1476793	17.24%	3.236%
40	14.269	1896	1901	1905	rVB	18495	24148	0.28%	0.053%
41	14.351	1908	1915	1921	rBV	1146163	1649457	19.26%	3.615%
42	14.540	1938	1947	1962	rBV	316997	487875	5.70%	1.069%
43	14.746	1978	1982	1983	rBV2	8748	11962	0.14%	0.026%
44	14.769	1983	1986	1991	rVB3	35726	44983	0.53%	0.099%
45	15.222	2059	2063	2067	rBV2	12283	16274	0.19%	0.036%
46	15.304	2073	2077	2078	rBV3	19113	19469	0.23%	0.043%
47	15.346	2078	2084	2091	rVB	1302791	1820685	21.26%	3.990%
48	15.457	2097	2103	2109	rBV	201187	285187	3.33%	0.625%
49	15.510	2109	2112	2119	rVB	28415	35021	0.41%	0.077%
50	15.587	2119	2125	2130	rBV3	11764	17255	0.20%	0.038%
51	15.775	2151	2157	2167	rBV	181545	270873	3.16%	0.594%
52	16.081	2205	2209	2213	rBV2	8411	11571	0.14%	0.025%
53	16.251	2234	2238	2242	rVB3	12019	16979	0.20%	0.037%
54	16.357	2251	2256	2260	rBV4	5826	9049	0.11%	0.020%
55	16.598	2293	2297	2302	rBV6	5105	9143	0.11%	0.020%
56	16.804	2328	2332	2335	rBV	7605	9217	0.11%	0.020%
57	16.875	2341	2344	2349	rBV3	21058	26963	0.31%	0.059%
58	16.928	2350	2353	2358	rVB4	5758	8651	0.10%	0.019%
59	17.098	2375	2382	2388	rBV	1423851	1995019	23.30%	4.372%
60	17.192	2392	2398	2406	rVV	1368219	2027779	23.68%	4.444%
61	17.292	2411	2415	2420	rVV8	19280	31789	0.37%	0.070%
62	17.416	2432	2436	2439	rVB6	5800	8735	0.10%	0.019%
63	17.475	2443	2446	2448	rBV4	7456	9301	0.11%	0.020%
64	17.616	2466	2470	2476	rBV2	31969	46383	0.54%	0.102%
65	17.787	2494	2499	2506	rVB	137566	186456	2.18%	0.409%
66	18.004	2532	2536	2543	rBV	92031	135951	1.59%	0.298%
67	18.251	2575	2578	2583	rVB2	18328	21271	0.25%	0.047%
68	18.310	2584	2588	2595	rVB	41238	62617	0.73%	0.137%
69	18.428	2603	2608	2614	rBV	289346	399002	4.66%	0.874%
70	18.963	2690	2699	2703	rBV2	425194	622435	7.27%	1.364%
71	19.098	2718	2722	2725	rBV	118779	126596	1.48%	0.277%
72	19.287	2752	2754	2757	rBV3	24181	27533	0.32%	0.060%
73	19.498	2784	2790	2799	rBV	1932430	2705734	31.59%	5.929%
74	20.063	2883	2886	2889	rVB2	39799	41782	0.49%	0.092%
75	20.216	2909	2912	2914	rBV4	17499	19802	0.23%	0.043%
76	20.298	2923	2926	2932	rVB6	23576	35889	0.42%	0.079%

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Misc :
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Instrument : BNA_N
ClientSampleId : BGDB6

Integration Parameters: LSCINT.P
Integrator: RTE
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Sampling : 1 Min Area: 0.1 % of largest Peak
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Title : SVOA CALIBRATION

77	20.386	2939	2941	2947	rVB	53998	58008	0.68%	0.127%
78	20.433	2947	2949	2958	rVB8	18789	19674	0.23%	0.043%
79	20.516	2958	2963	2969	rBV	7274259	8563821	100.00%	18.766%
80	20.675	2987	2990	2993	rBV5	21009	22704	0.27%	0.050%
81	20.716	2994	2997	3003	rBV6	64401	96283	1.12%	0.211%
82	20.910	3026	3030	3034	rBV2	123962	144980	1.69%	0.318%
83	21.028	3046	3050	3051	rBV	134778	174939	2.04%	0.383%
84	21.051	3051	3054	3065	rVB	421438	540562	6.31%	1.185%
85	21.222	3079	3083	3088	rBV	290903	316102	3.69%	0.693%
86	21.292	3090	3095	3101	rBV	1722469	2209869	25.80%	4.843%
87	21.463	3122	3124	3136	rVB4	50468	82612	0.96%	0.181%
88	21.904	3195	3199	3207	rBV2	80412	150412	1.76%	0.330%
89	22.157	3238	3242	3248	rBV	203336	251609	2.94%	0.551%
90	22.369	3275	3278	3287	rVB	92462	129164	1.51%	0.283%
91	22.792	3346	3350	3357	rVB3	153142	224280	2.62%	0.491%
92	22.875	3360	3364	3371	rBV	155770	229425	2.68%	0.503%
93	23.398	3446	3453	3459	rBV	1380810	2635152	30.77%	5.775%
94	23.451	3459	3462	3469	rVB2	162605	245177	2.86%	0.537%
95	23.539	3470	3477	3485	rBV2	1223153	2318336	27.07%	5.080%
96	24.098	3567	3572	3579	rVB	165934	316034	3.69%	0.693%
97	24.851	3695	3700	3709	rVB3	146031	328010	3.83%	0.719%

Sum of corrected areas: 45633581

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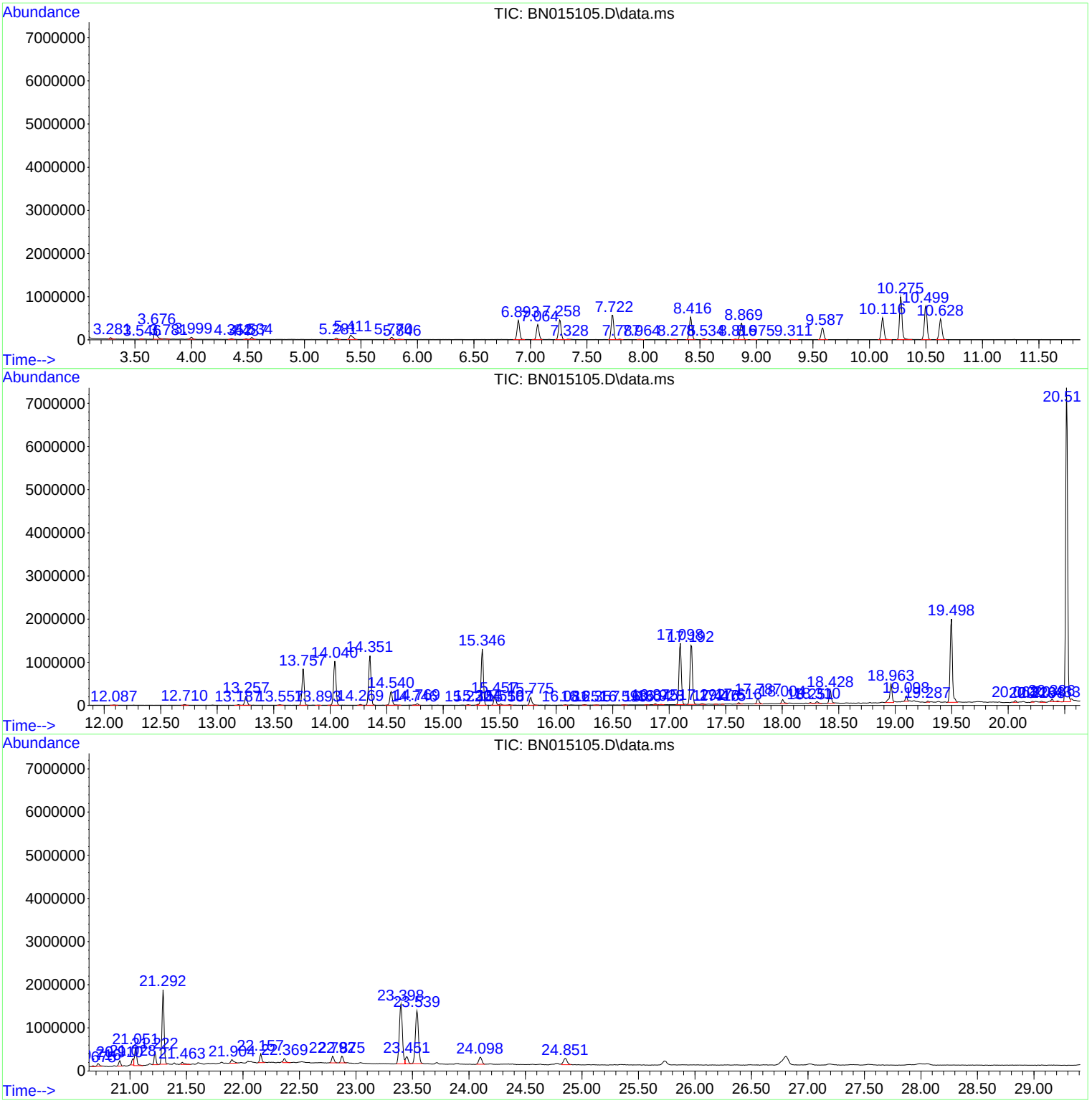
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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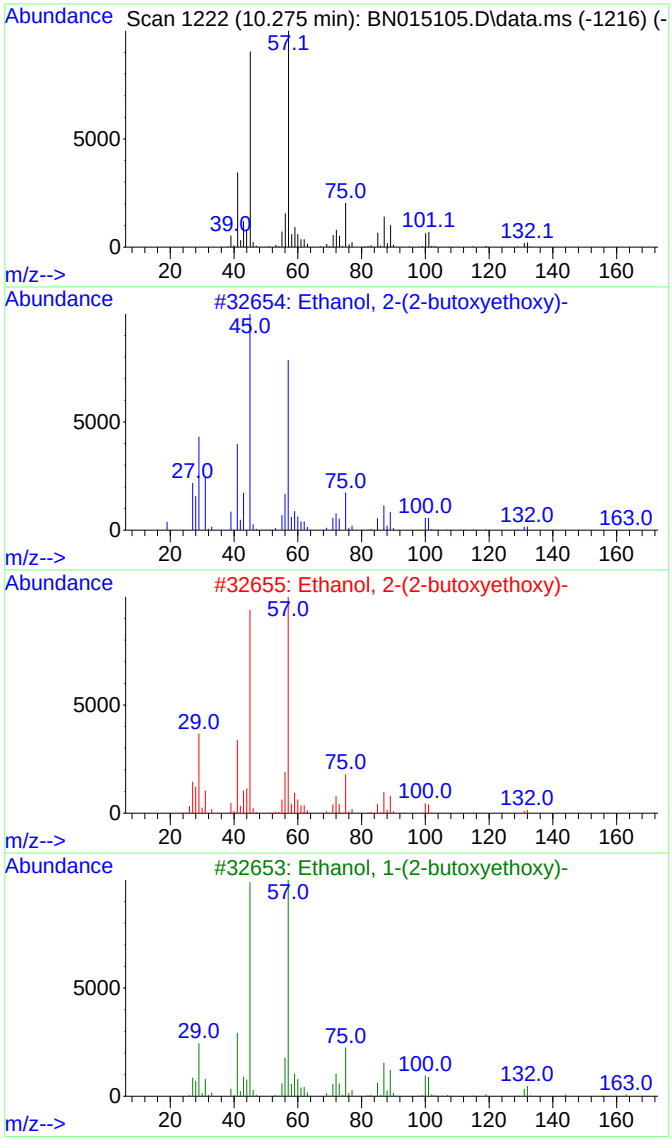
Quant Title : SVOA CALIBRATION

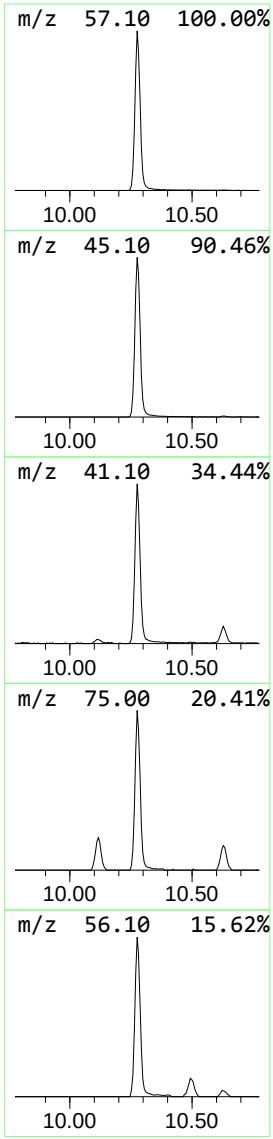
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	23.89 ng/ul	1526750	Naphthalene-d8	10.499

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-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





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Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

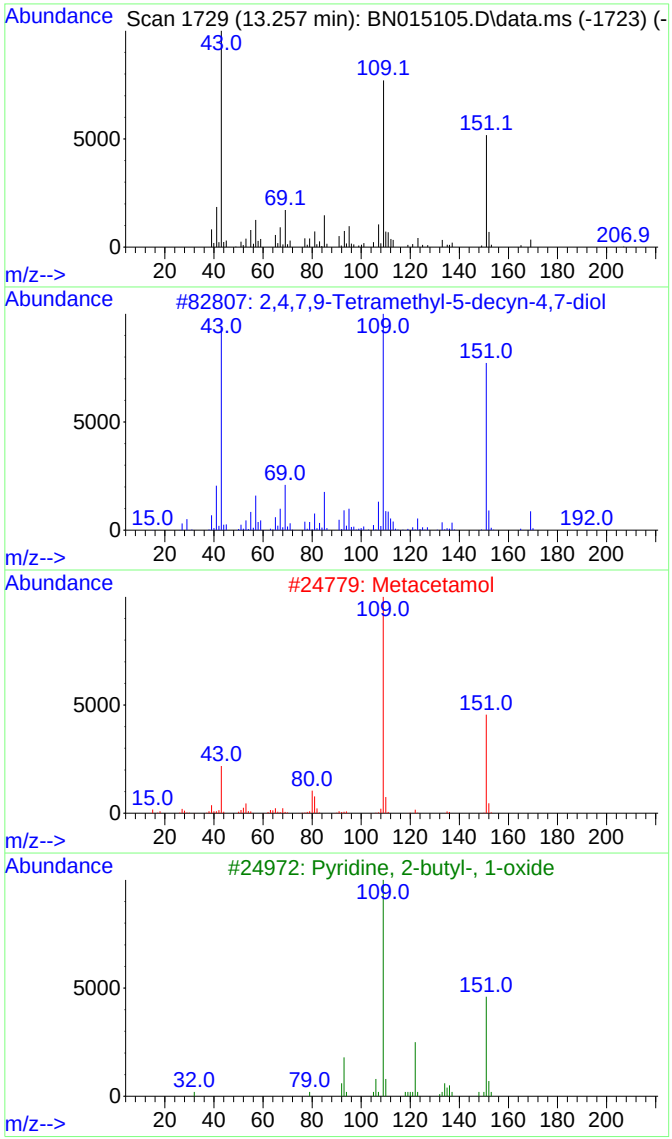
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

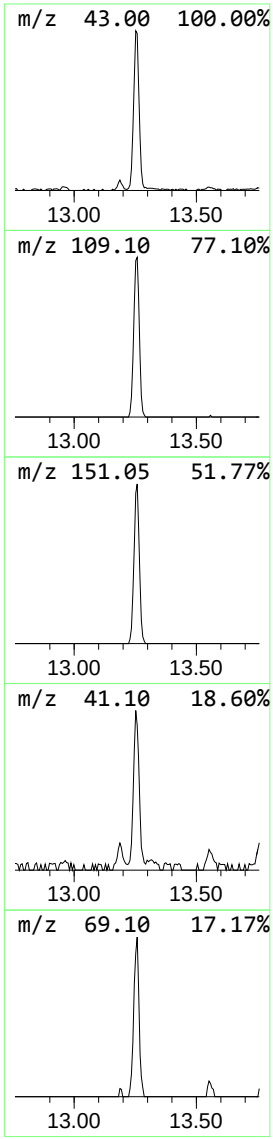
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.257	3.83 ng/ul	315526	Acenaphthene-d10	14.352

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	Metacetamol	151	C8H9NO2	000621-42-1	53
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53
4	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	53
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53





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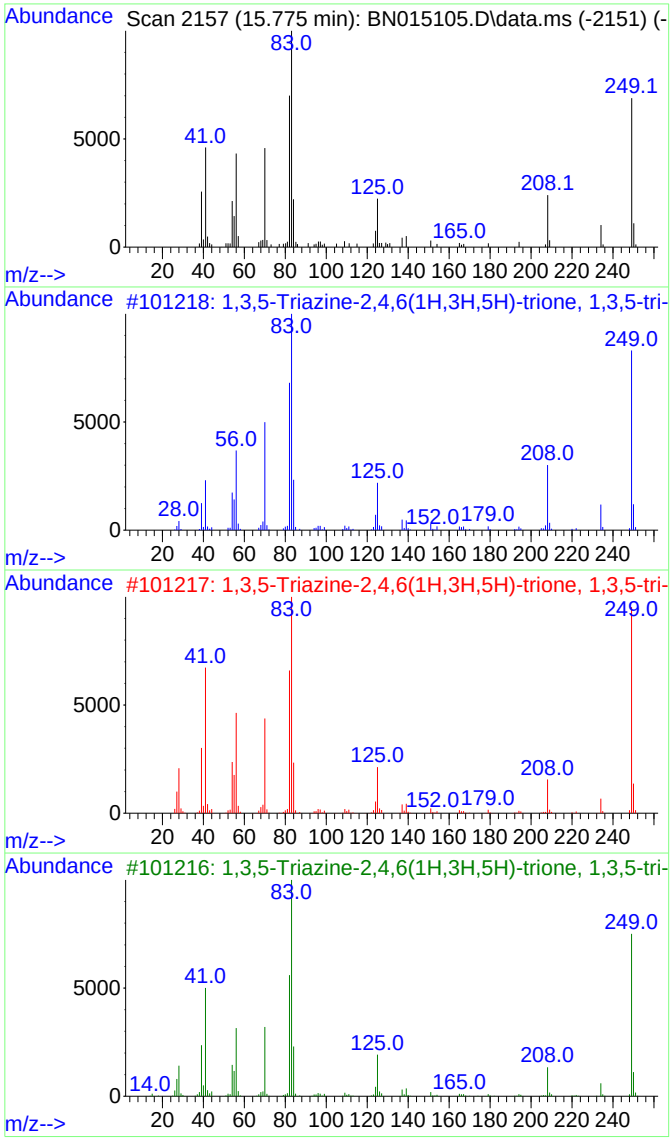
Quant Title : SVOA CALIBRATION

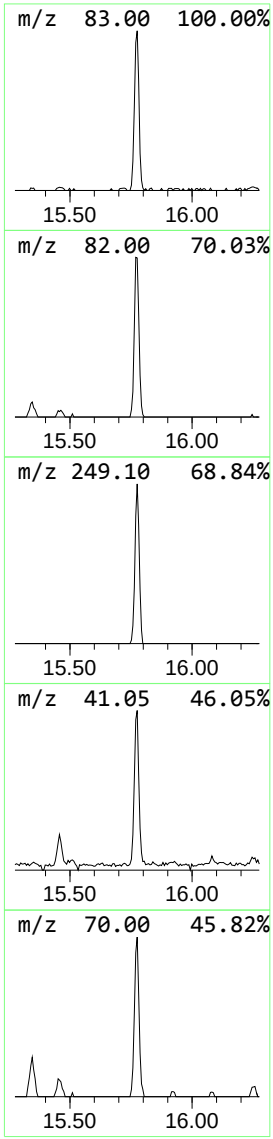
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.72 ng/ul	270873	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
2	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
3	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	93
4	Azetidine, 2-methyl-1-[3,3,3-tri...	249	C8H9F6NO	050837-79-1	35
5	N-Carboxy-1-[4-chlorophenyl]-2-[...	295	C15H18ClNO3	058158-38-6	23





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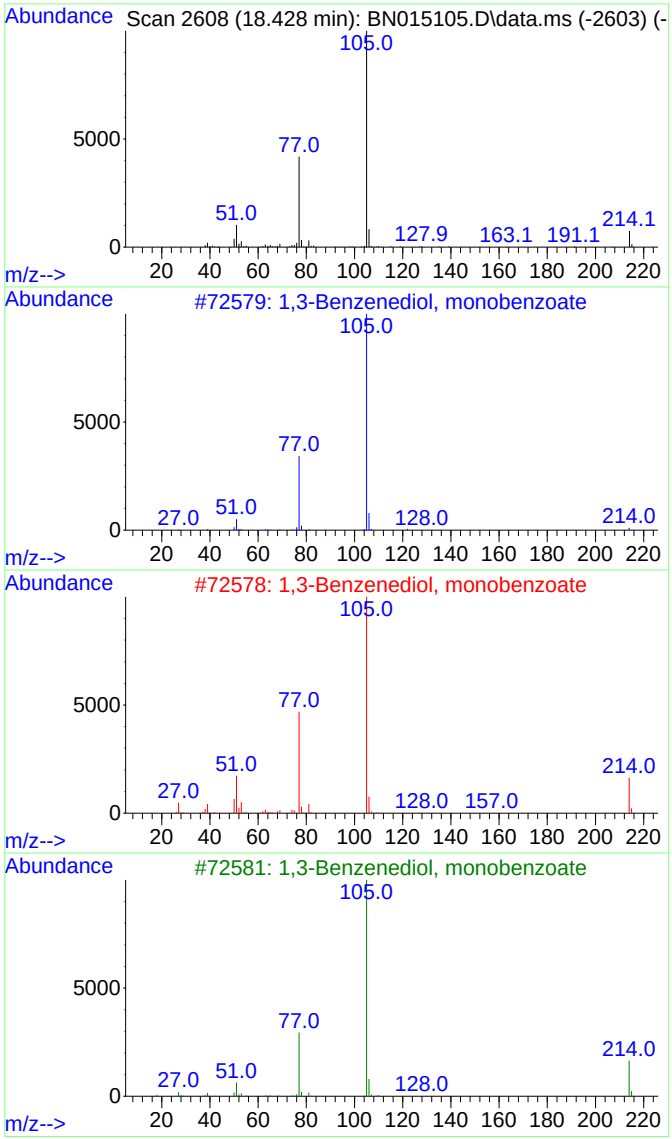
Quant Title : SVOA CALIBRATION

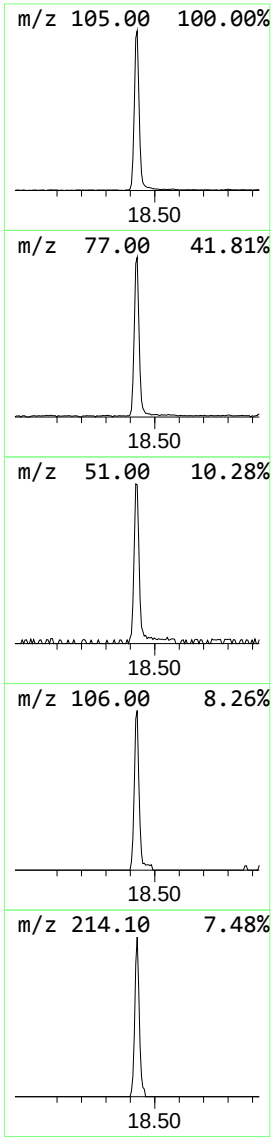
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,3-Benzenediol, monobenzoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.428	4.00 ng/ul	399002	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	91
2	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90
3	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	87
4	1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	78
5	Isopropyl phenyl ketone	148	C10H12O	000611-70-1	72





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Instrument : BNA_N
ClientSampleId : BGDB6

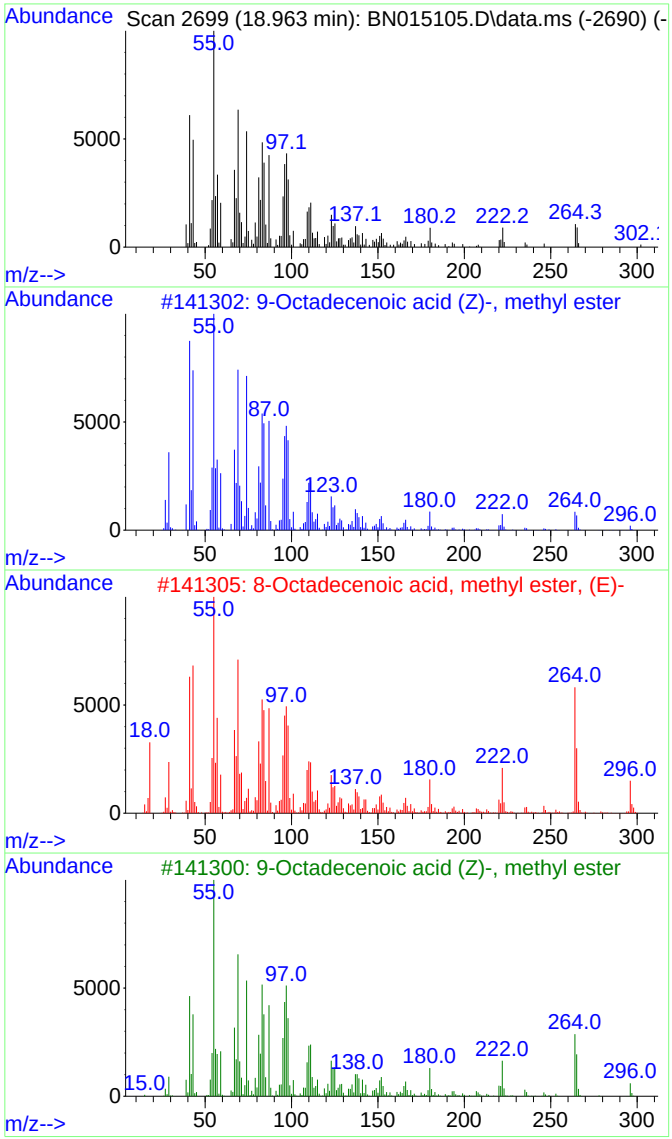
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

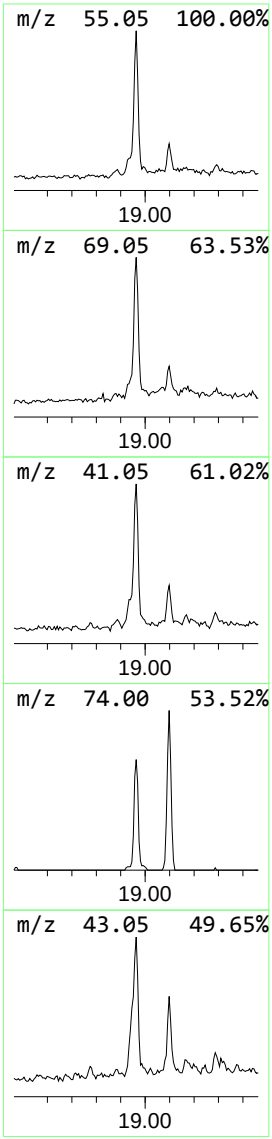
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.963	6.24 ng/ul	622435	Phenanthrene-d10	17.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	95
2	8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	94
3	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	94
4	8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	94
5	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94





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Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB6

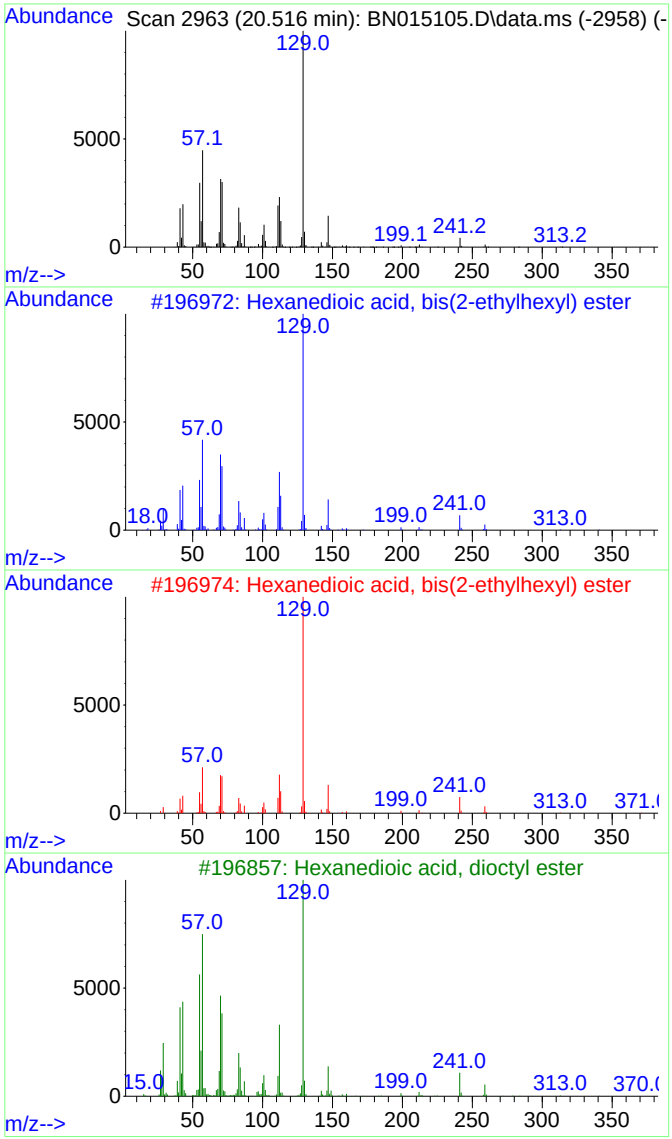
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

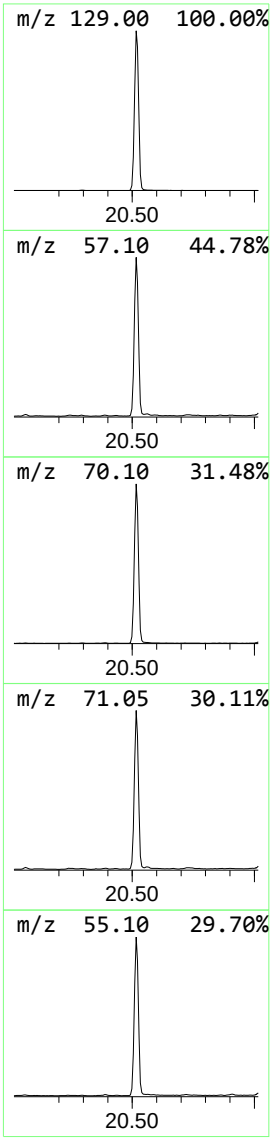
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	77.51 ng/ul	8563820	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76





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Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

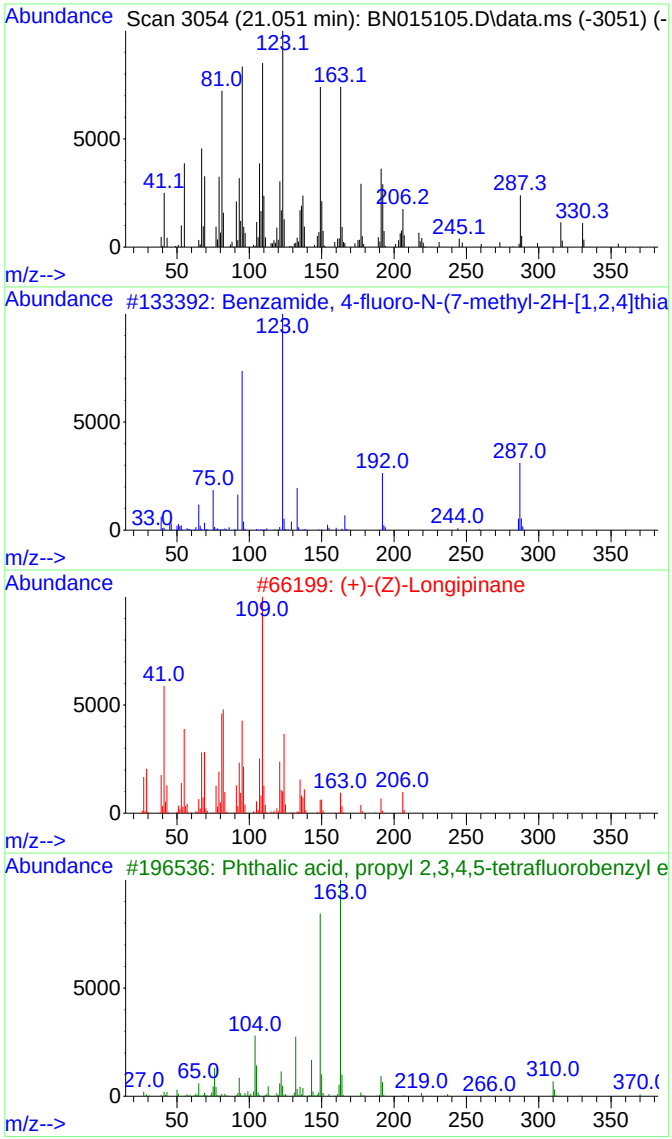
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

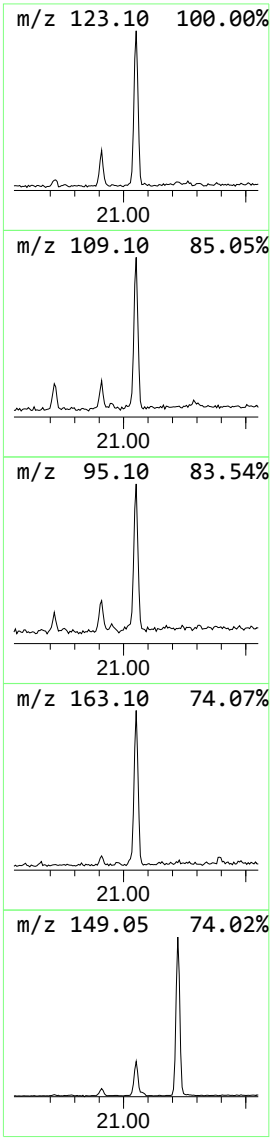
****

Peak Number 8 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.051	4.89 ng/ul	540562	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, 4-fluoro-N-(7-methyl-...	287	C14H10FN3OS	1000350-88-7	27
2	(+)-(Z)-Longipinane	206	C15H26	1000156-12-3	25
3	Phthalic acid, propyl 2,3,4,5-te...	370	C18H14F4O4	1000377-72-5	14
4	3-Fluorobenzoic acid, hexyl ester	224	C13H17FO2	1000279-00-8	14
5	Hexahydroindole	123	C8H13N	018159-32-5	14





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Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

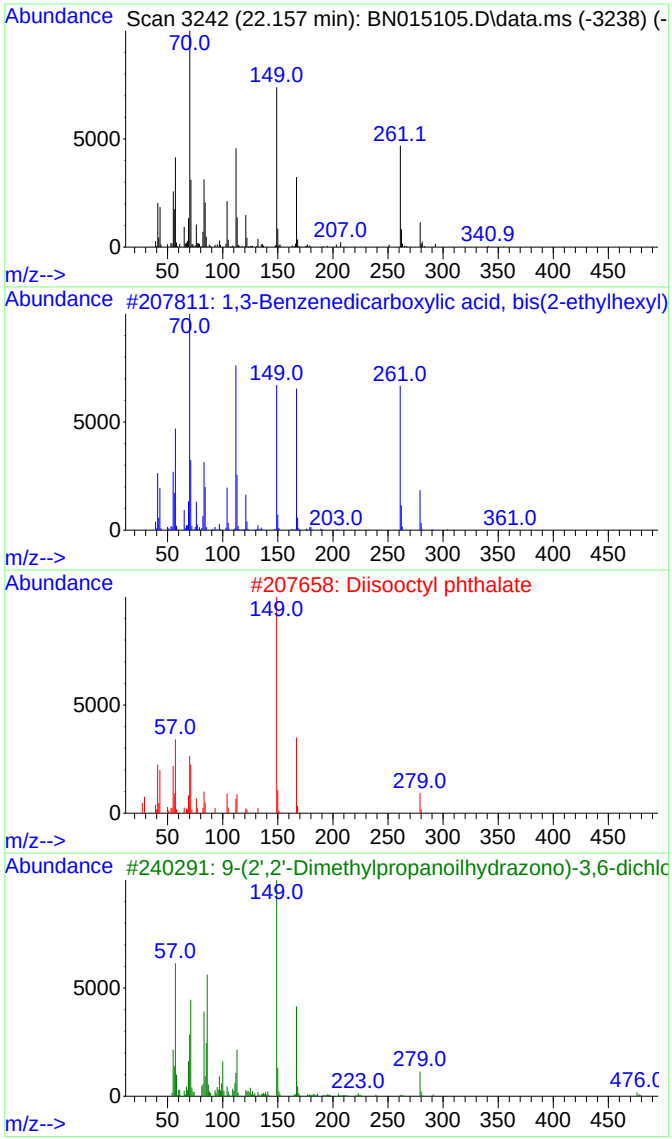
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

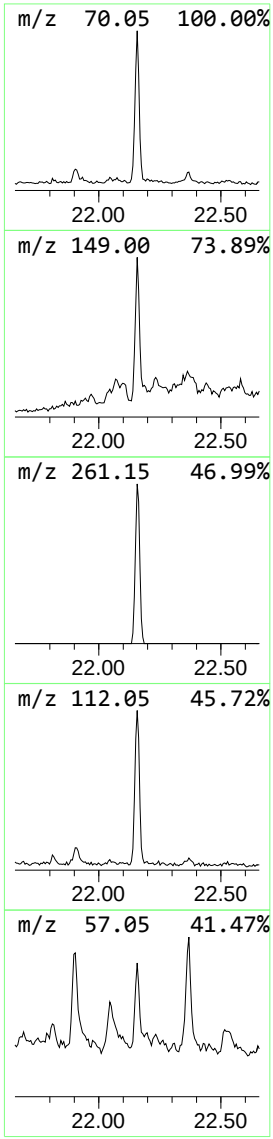
****

Peak Number 9 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.157	2.28 ng/ul	251609	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	49
2	Diisooctyl phthalate	390	C24H38O4	000131-20-4	35
3	9-(2',2'-Dimethylpropanoilhydraz...	576	C30H42Cl2N4O3	1000111-04-6	27
4	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	27
5	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	27





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Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

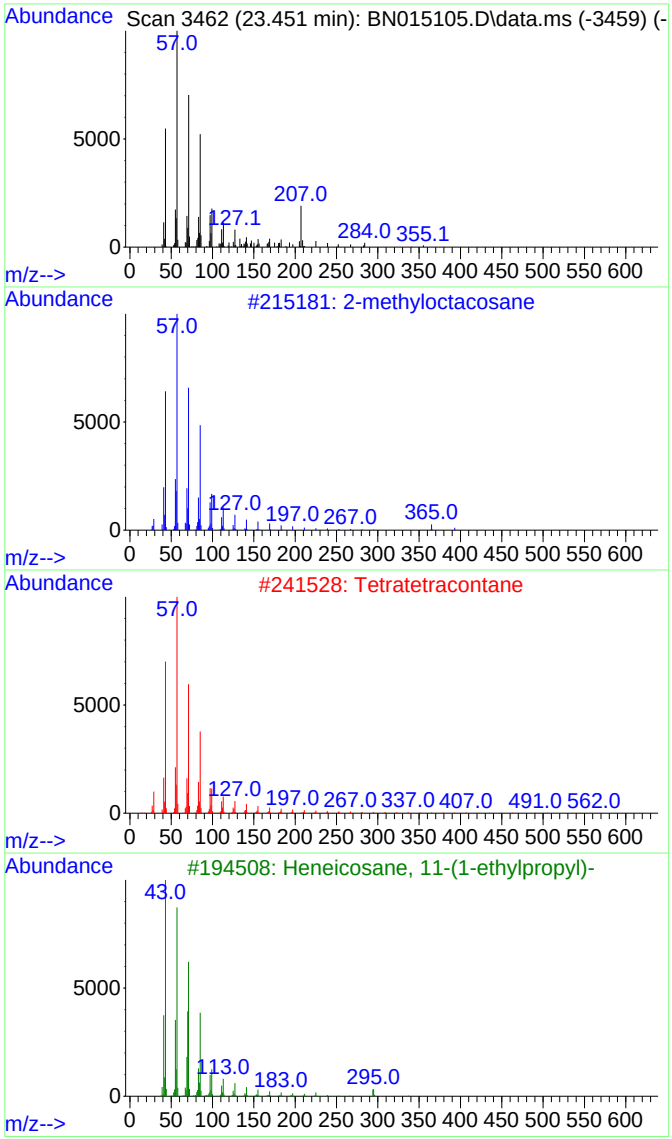
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

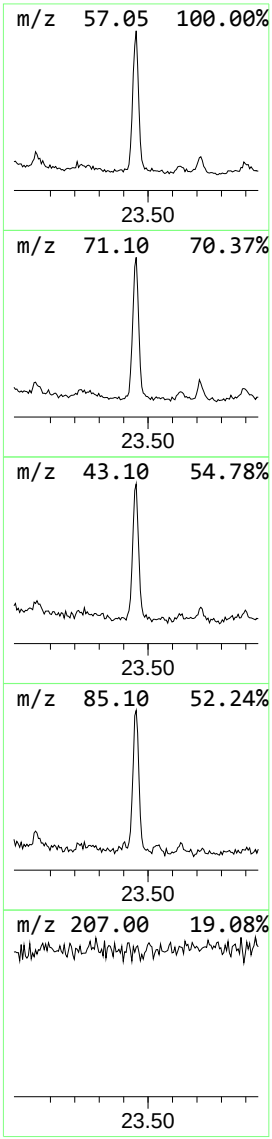
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.451	2.12 ng/ul	245177	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane	408	C29H60	1000376-72-8	87
2	Tetratetracontane	619	C44H90	007098-22-8	83
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83
5	Heneicosane	296	C21H44	000629-94-7	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

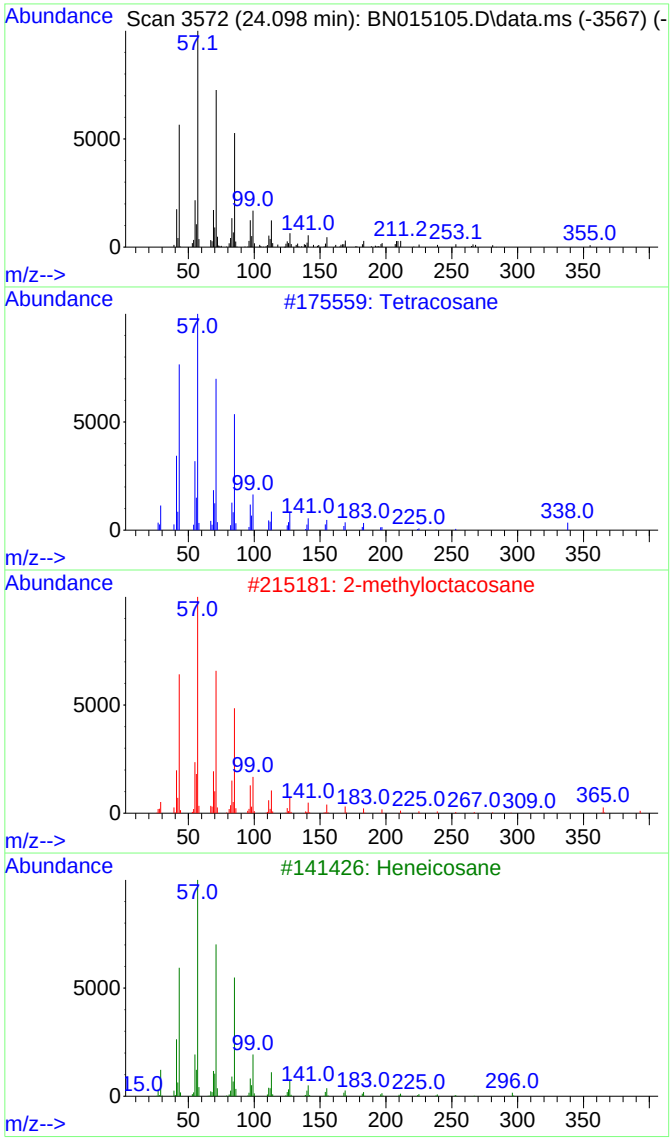
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

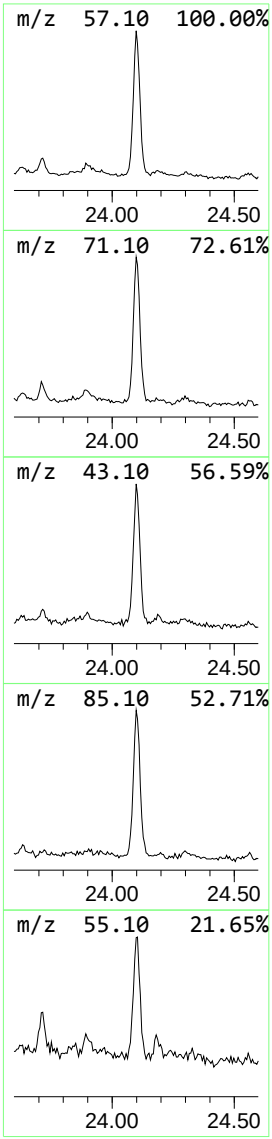
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.098	2.73 ng/ul	316034	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	2-methyloctacosane	408	C29H60	1000376-72-8	90
3	Heneicosane	296	C21H44	000629-94-7	90
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5	Docosane	310	C22H46	000629-97-0	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDB6

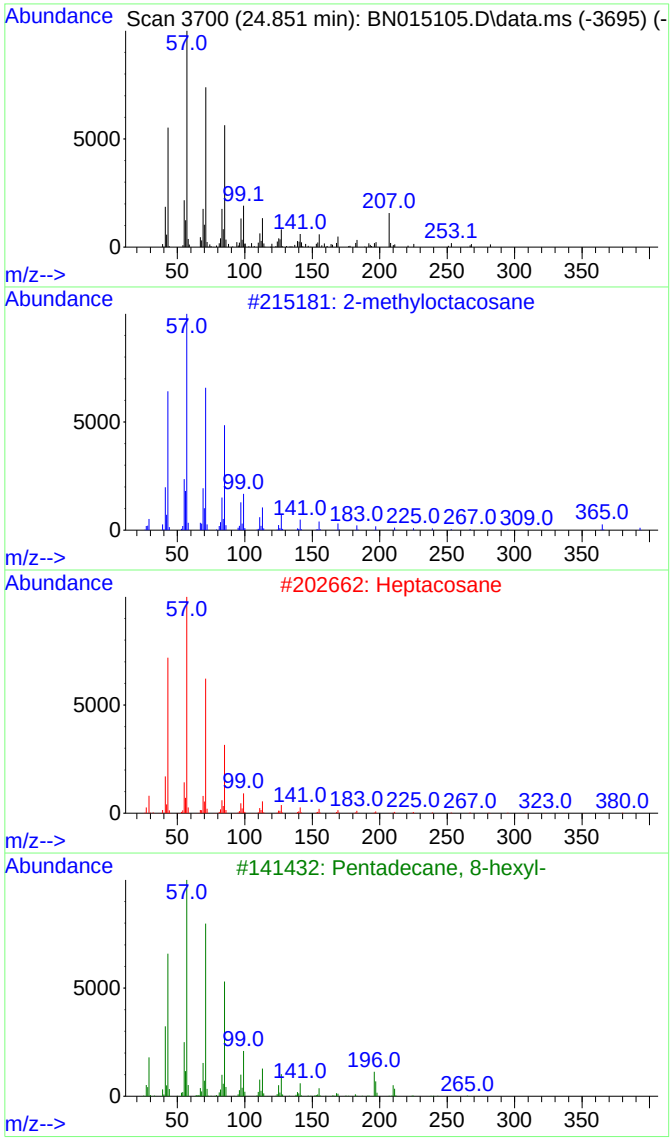
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

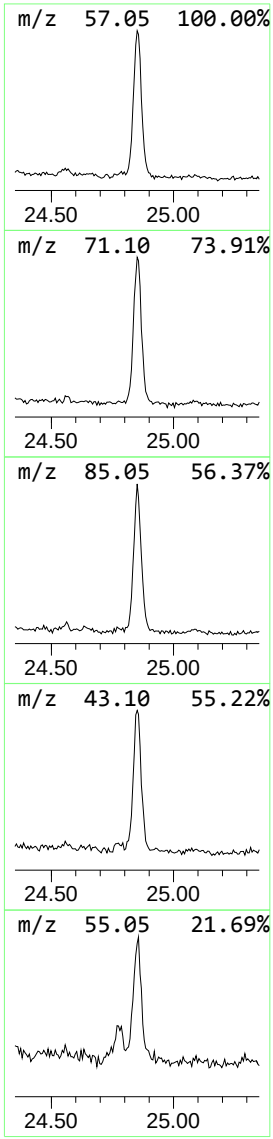
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.851	2.83 ng/ul	328010	Perylene-d12	23.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane	408	C29H60	1000376-72-8	90
2	Heptacosane	380	C27H56	000593-49-7	87
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87
4	Nonadecane	268	C19H40	000629-92-5	87
5	Eicosane, 2-methyl-	296	C21H44	001560-84-5	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015105.D
Acq On : 26 Jun 2021 08:02
Operator : CG/JU
Sample : M2704-07
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BGDB6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.275	23.9	ng/ul	1526750	2	10.499	1278080	20.0
2,4,7,9-Tetrame...	13.257	3.8	ng/ul	315526	3	14.352	1649460	20.0
1,3,5-Triazine-...	15.775	2.7	ng/ul	270873	4	17.098	1995020	20.0
1,3-Benzenediol...	18.428	4.0	ng/ul	399002	4	17.098	1995020	20.0
9-Octadecenoic ...	18.963	6.2	ng/ul	622435	4	17.098	1995020	20.0
Hexanedioic aci...	20.516	77.5	ng/ul	8563820	5	21.292	2209870	20.0
unknown-01	21.051	4.9	ng/ul	540562	5	21.292	2209870	20.0
unknown-02	22.157	2.3	ng/ul	251609	5	21.292	2209870	20.0
(DEL) Alkane: S...	23.451	2.1	ng/ul	245177	6	23.539	2318340	20.0
(DEL) Alkane: S...	24.098	2.7	ng/ul	316034	6	23.539	2318340	20.0
(DEL) Alkane: S...	24.851	2.8	ng/ul	328010	6	23.539	2318340	20.0