LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

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-BP072421.M
Title : SVOA CALIBRATION

Signal : TIC: BP006353.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.305	33	37	46	rVB	71383	102928	1.52%	0.140%
2	3.711	103	106	125	rBV	206477	374159	5.54%	0.510%
3	3.934	138	144	152	rBV3	20616	38336	0.57%	0.052%
4	4.023	155	159	166	rVB	22828	35105	0.52%	0.048%
5	4.381	217	220	226	rVB	29702	39037	0.58%	0.053%
6	4.523	241	244	249	rVB	11407	16258	0.24%	0.022%
7	5.317	374	379	383	rVB	15401	22361	0.33%	0.030%
8	5.858	466	471	474	rVV	111926	169630	2.51%	0.231%
9	5.893	474	477	486	rVB	125652	181053	2.68%	0.247%
10	6.452	567	572	578	rVB2	41314	62924	0.93%	0.086%
11	6.905	644	649	650	rBV4	11865	14734	0.22%	0.020%
12	6.958	650	658	671	rVB2	266399	502843	7.44%	0.685%
13	7.128	681	687	698	rVB	1028295	1640419	24.27%	2.234%
14	7.317	712	719	731	rBV	941024	1501544	22.22%	2.045%
15	7.417	731	736	740	rVB6	8918	14835	0.22%	0.020%
16	7.593	755	766	773	rVB4	9311	27070	0.40%	0.037%
17	7.787	791	799	805	rBV	1018126	1586662	23.48%	2.161%
18	7.858	805	811	821	rVB	145559	243310	3.60%	0.331%
19	7.946	821	826	833	rVB9	7771	16647	0.25%	0.023%
20	8.028	833	840	847	rBV2	41211	79636	1.18%	0.108%
21	8.093	847	851	856	rVB8	8793	16516	0.24%	0.022%
22	8.411	898	905	908	rVB4	17386	30617	0.45%	0.042%
23	8.493	912	919	928	rVV	738498	1181531	17.48%	1.609%
24	8.581	928	934	936	rVV	42583	69533	1.03%	0.095%
25	8.611	936	939	945	rVB	46436	68941	1.02%	0.094%
26	8.881	978	985	989	rBV	303584	478894	7.09%	0.652%
27	8.940	989	995	1002	rVV	1231565	1980122	29.30%	2.697%
28	9.034	1006	1011	1019	rVV	92688	146927	2.17%	0.200%
29	9.116	1019	1025	1031	rVB	115799	179093	2.65%	0.244%
30	9.505	1085	1091	1098	rVB	8492	20803	0.31%	0.028%
31	9.658	1110	1117	1128	rBV	611124	1007019	14.90%	1.372%
32	9.746	1128	1132	1137	rVV	20551	33426	0.49%	0.046%
33	9.840	1138	1148	1154	rVV	553968	1265193	18.72%	1.723%
34	9.887	1154	1156	1164	rVB9	14764	25338	0.37%	0.035%
35	10.081	1178	1189	1197	rBV2	797340	1705459	25.23%	2.323%
36	10.193	1201	1208	1218	rVB	1242186	2017962	29.86%	2.748%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
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Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

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-BP072421.M
Title : SVOA CALIBRATION

37	10.358	1230	1236	1242	rVB4	23576	42044	0.62%	0.057%
38	10.440	1244	1250	1254	rBV5	30349	48494	0.72%	0.066%
39	10.487	1254	1258	1264	rVV4	46035	79232	1.17%	0.108%
40	10.569	1264	1272	1278	rVV	1350356	2245028	33.22%	3.058%
41	10.634	1278	1283	1289	rVV3	59611	114314	1.69%	0.156%
42	10.710	1289	1296	1307	rVB	1377520	2255265	33.37%	3.072%
43	10.881	1319	1325	1330	rBV7	15519	29448	0.44%	0.040%
44	10.957	1330	1338	1347	rVB7	8832	29687	0.44%	0.040%
45	11.122	1360	1366	1371	rVV	128690	200664	2.97%	0.273%
46	11.181	1371	1376	1380	rVV	140795	217852	3.22%	0.297%
47	11.222	1380	1383	1394	rVB2	48167	84170	1.25%	0.115%
48	11.322	1394	1400	1403	rBV5	10790	16589	0.25%	0.023%
49	11.369	1403	1408	1416	rVV5	21621	40182	0.59%	0.055%
50	11.746	1467	1472	1476	rBV3	16764	25754	0.38%	0.035%
51	11.805	1478	1482	1487	rVB3	14123	19703	0.29%	0.027%
52	11.869	1489	1493	1494	rVV	14332	16716	0.25%	0.023%
53	11.893	1494	1497	1505	rVB3	16809	33520	0.50%	0.046%
54	11.999	1510	1515	1523	rVB6	17537	33710	0.50%	0.046%
55	12.157	1536	1542	1548	rBV	27850	48688	0.72%	0.066%
56	12.493	1595	1599	1604	rVV6	12462	20801	0.31%	0.028%
57	12.775	1640	1647	1654	rBV	133910	208553	3.09%	0.284%
58	13.028	1685	1690	1696	rVV	220283	318645	4.71%	0.434%
59	13.263	1725	1730	1734	rBV2	13999	20035	0.30%	0.027%
60	13.334	1736	1742	1748	rVV	221606	349078	5.17%	0.475%
61	13.381	1748	1750	1755	rVB4	10594	15860	0.23%	0.022%
62	13.657	1790	1797	1807	rBV3	59739	109379	1.62%	0.149%
63	13.834	1820	1827	1837	rBV	2627953	3608166	53.39%	4.914%
64	13.951	1842	1847	1854	rVB6	23513	39508	0.58%	0.054%
65	14.104	1861	1873	1880	rBV	2909280	4151638	61.43%	5.654%
66	14.275	1899	1902	1907	rVV5	9787	16267	0.24%	0.022%
67	14.410	1915	1925	1934	rBV	1863123	2822508	41.76%	3.844%
68	14.504	1936	1941	1946	rVB	20972	29933	0.44%	0.041%
69	14.610	1953	1959	1967	rBV	102538	165728	2.45%	0.226%
70	14.675	1968	1970	1975	rVB6	10062	14643	0.22%	0.020%
71	14.957	2013	2018	2023	rBV3	14989	23001	0.34%	0.031%
72	15.016	2023	2028	2033	rBV3	27738	41711	0.62%	0.057%
73	15.163	2048	2053	2059	rVB	192983	250037	3.70%	0.341%
74	15.228	2059	2064	2066	rBV	45557	63942	0.95%	0.087%
75	15.410	2087	2095	2107	rBV	3471207	5197698	76.91%	7.079%
76	15.522	2109	2114	2117	rVV3	22378	36775	0.54%	0.050%

LSC Area Percent Report

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Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

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-BP072421.M
Title : SVOA CALIBRATION

77	15.557	2117	2120	2125	rVV3	17662	26805	0.40%	0.037%
78	15.645	2129	2135	2143	rBV2	48073	76732	1.14%	0.105%
79	15.781	2152	2158	2165	rBV2	38426	58980	0.87%	0.080%
80	16.675	2303	2310	2314	rBV3	22438	36862	0.55%	0.050%
81	16.769	2322	2326	2331	rBV4	10058	15878	0.23%	0.022%
82	16.945	2352	2356	2362	rVB4	11087	19030	0.28%	0.026%
83	17.163	2386	2393	2400	rVB	2246581	3113350	46.07%	4.240%
84	17.263	2403	2410	2420	rBV	4346844	6030756	89.23%	8.214%
85	17.469	2441	2445	2453	rVB	19546	29934	0.44%	0.041%
86	17.545	2453	2458	2462	rBV7	10094	19924	0.29%	0.027%
87	17.845	2503	2509	2519	rBV	1912566	2250902	33.31%	3.066%
88	18.063	2542	2546	2555	rBV3	37847	60568	0.90%	0.082%
89	18.139	2555	2559	2565	rBV2	42378	80130	1.19%	0.109%
90	18.310	2584	2588	2596	rVB5	18281	27321	0.40%	0.037%
91	19.081	2715	2719	2724	rVB2	50219	62024	0.92%	0.084%
92	19.145	2726	2730	2734	rBV	50458	67010	0.99%	0.091%
93	19.304	2754	2757	2762	rVB6	23788	29876	0.44%	0.041%
94	19.445	2778	2781	2785	rBV2	32802	37207	0.55%	0.051%
95	19.504	2785	2791	2797	rBV	5269565	6636133	98.19%	9.038%
96	19.551	2797	2799	2806	rVB	56012	79539	1.18%	0.108%
97	21.257	3084	3089	3096	rBV	2698809	3324989	49.20%	4.529%
98	22.369	3273	3278	3290	rBV2	722106	1154875	17.09%	1.573%
99	23.457	3455	3463	3472	rBV2	3492388	6758347	100.00%	9.205%
100	23.604	3481	3488	3498	rVB2	1821320	3444049	50.96%	4.691%

Sum of corrected areas: 73423052

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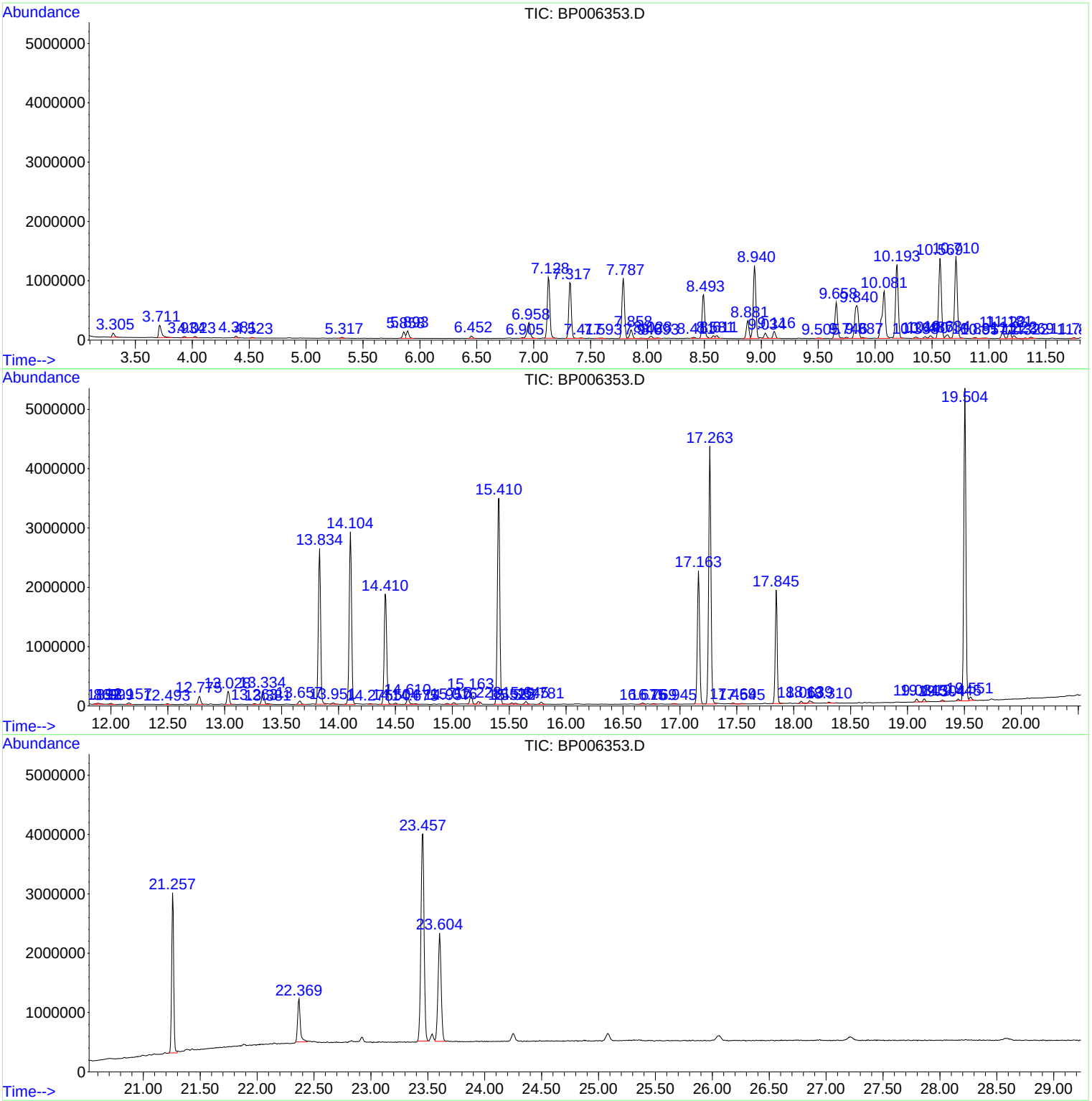
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

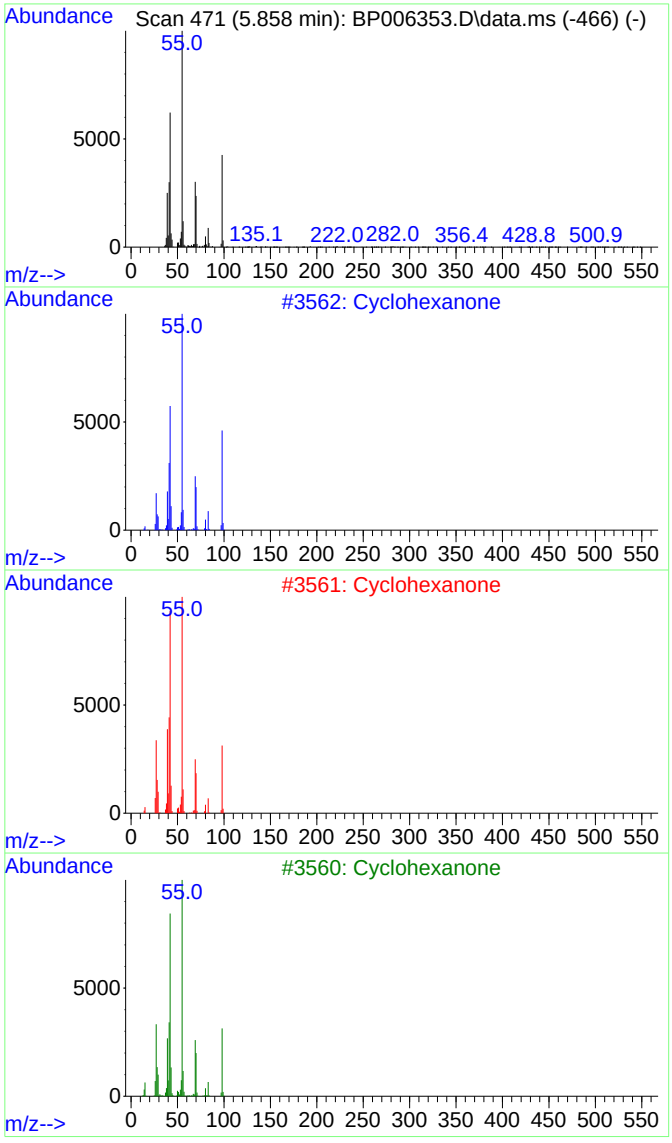
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

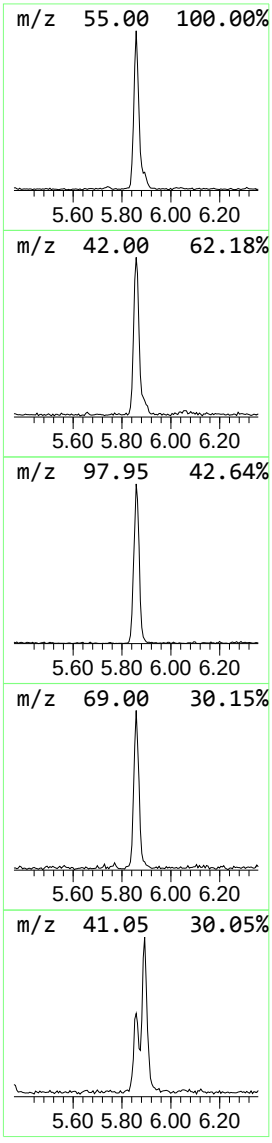
****

Peak Number 1 Cyclohexanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.860	2.14 ng/ul	169630	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone	98	C6H10O	000108-94-1	91
2	Cyclohexanone	98	C6H10O	000108-94-1	91
3	Cyclohexanone	98	C6H10O	000108-94-1	91
4	Cyclohexanone	98	C6H10O	000108-94-1	83
5	Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

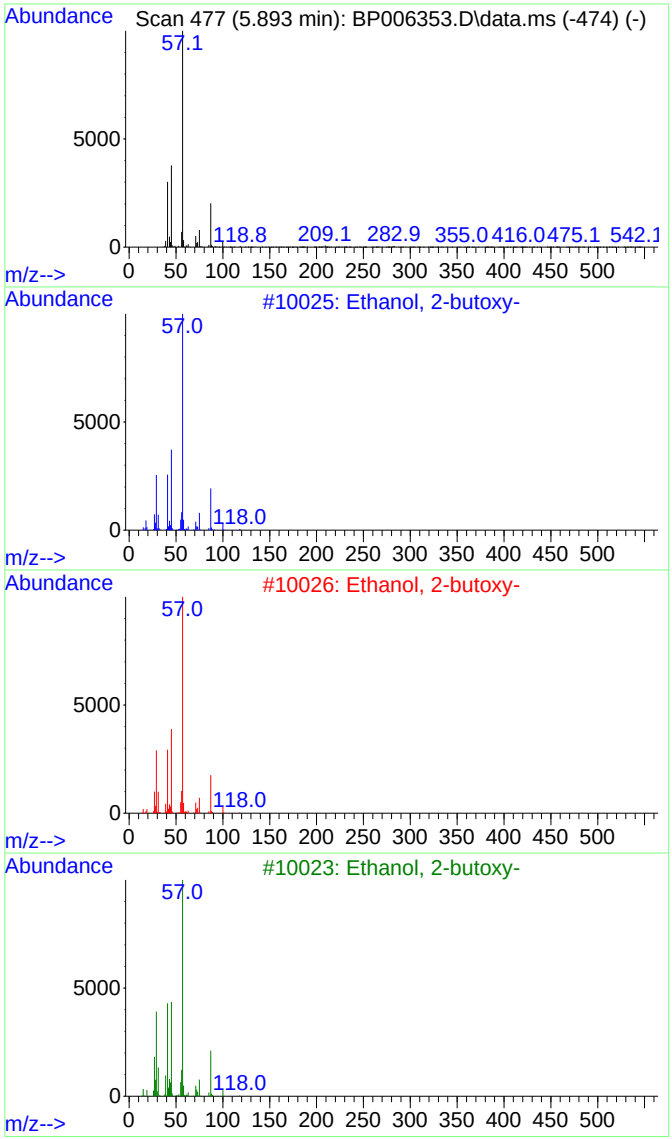
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

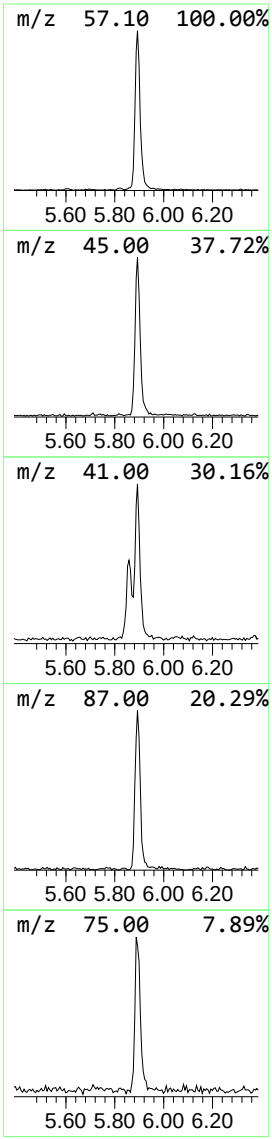
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Ethanol, 2-butoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.890	2.28 ng/ul	181053	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	72
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	53
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	50
4	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	47
5	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	38





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Acq On : 26 Jul 2021 23:44
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Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

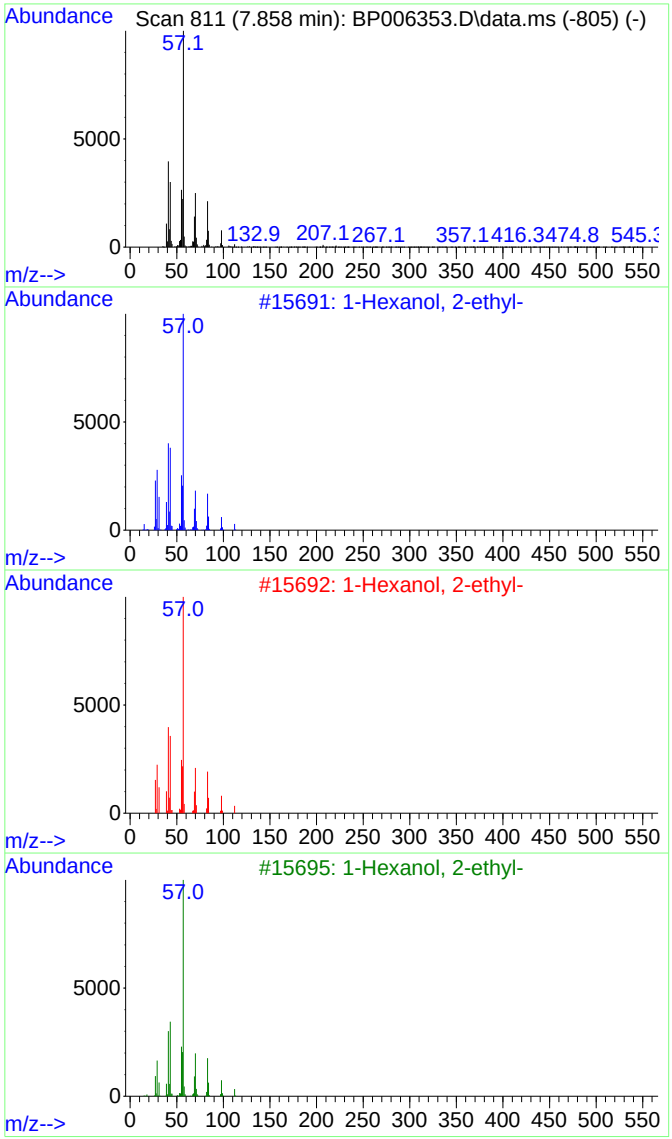
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

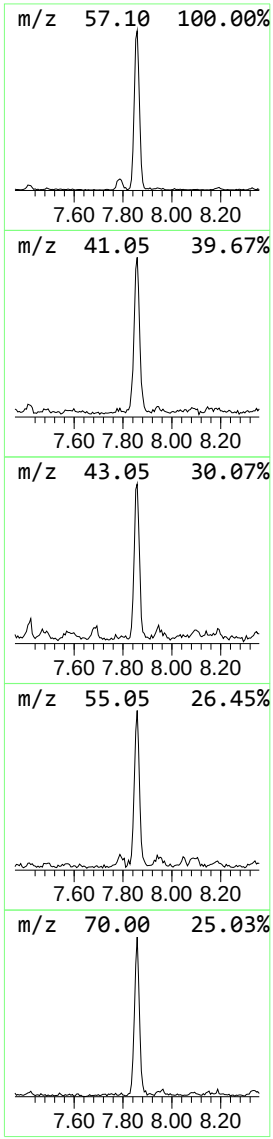
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.860	3.07 ng/ul	243310	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
5	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	59





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Data File : BP006353.D
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Instrument : BNA_P
ClientSampleId : C0060

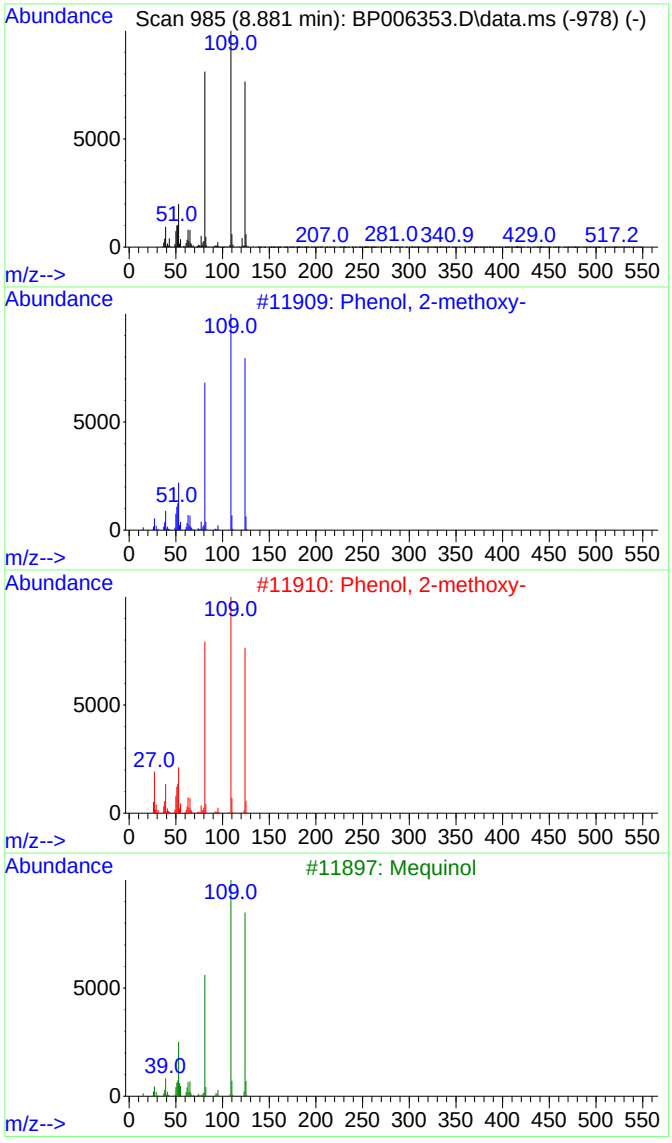
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

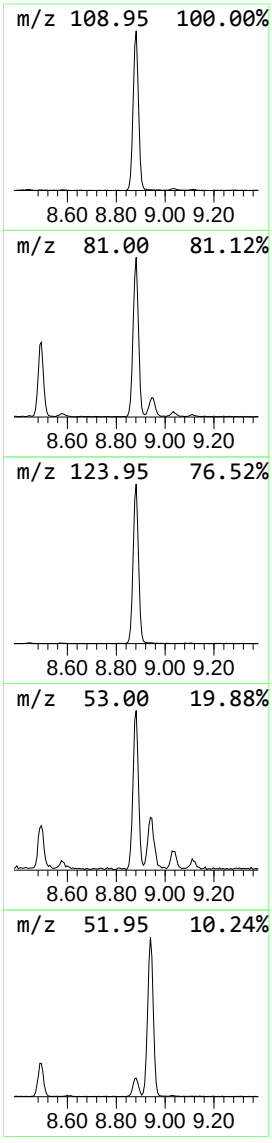
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 2-methoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.880	6.04 ng/ul	478894	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	96
2	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	94
3	Mequinol	124	C7H8O2	000150-76-5	91
4	Mequinol	124	C7H8O2	000150-76-5	87
5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

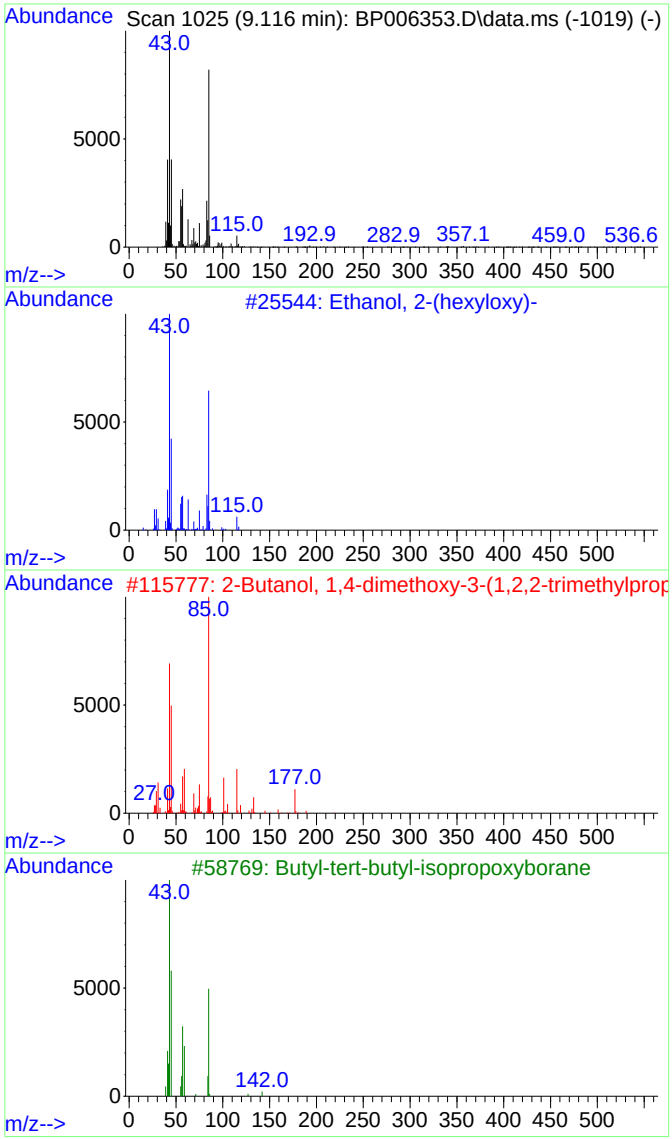
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

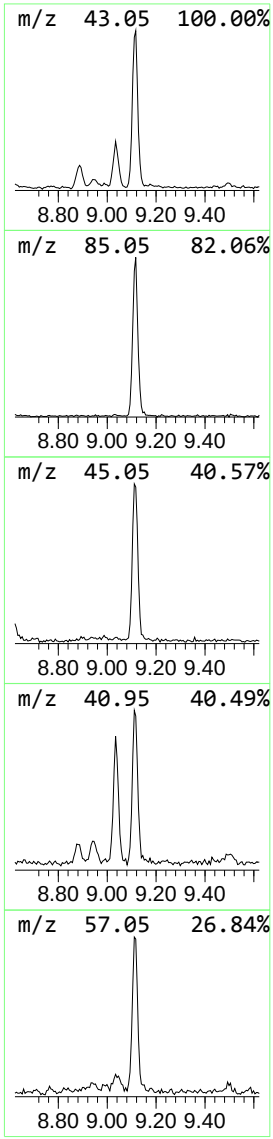
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-(hexyloxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.120	2.26 ng/ul	179093	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	86
2	2-Butanol, 1,4-dimethoxy-3-(1,2,...	234	C12H26O4	079449-92-6	59
3	Butyl-tert-butyl-isopropoxyborane	184	C11H25BO	097782-82-6	53
4	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	53
5	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	53





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Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

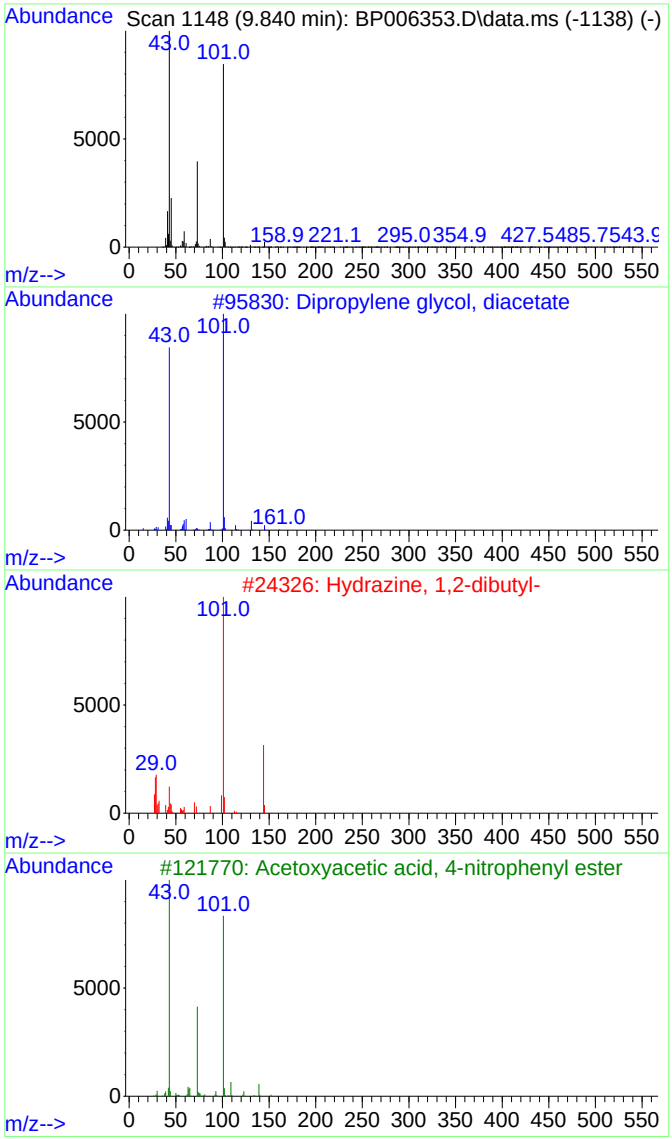
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

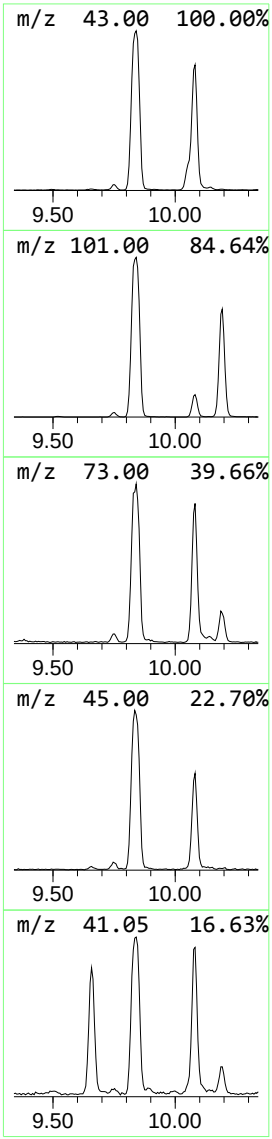
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Dipropylene glycol, diacetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.840	11.27 ng/ul	1265190	Naphthalene-d8	10.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dipropylene glycol, diacetate	218	C10H18O5	1000506-25-8	59
2	Hydrazine, 1,2-dibutyl-	144	C8H20N2	001744-71-4	49
3	Acetoxyacetic acid, 4-nitropheny...	239	C10H9NO6	1000307-54-5	45
4	4-Ketopimelic	174	C7H10O5	000502-50-1	45
5	Methyl(methyl 2,3,4-tri-O-methyl...	264	C11H20O7	052729-97-2	43





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Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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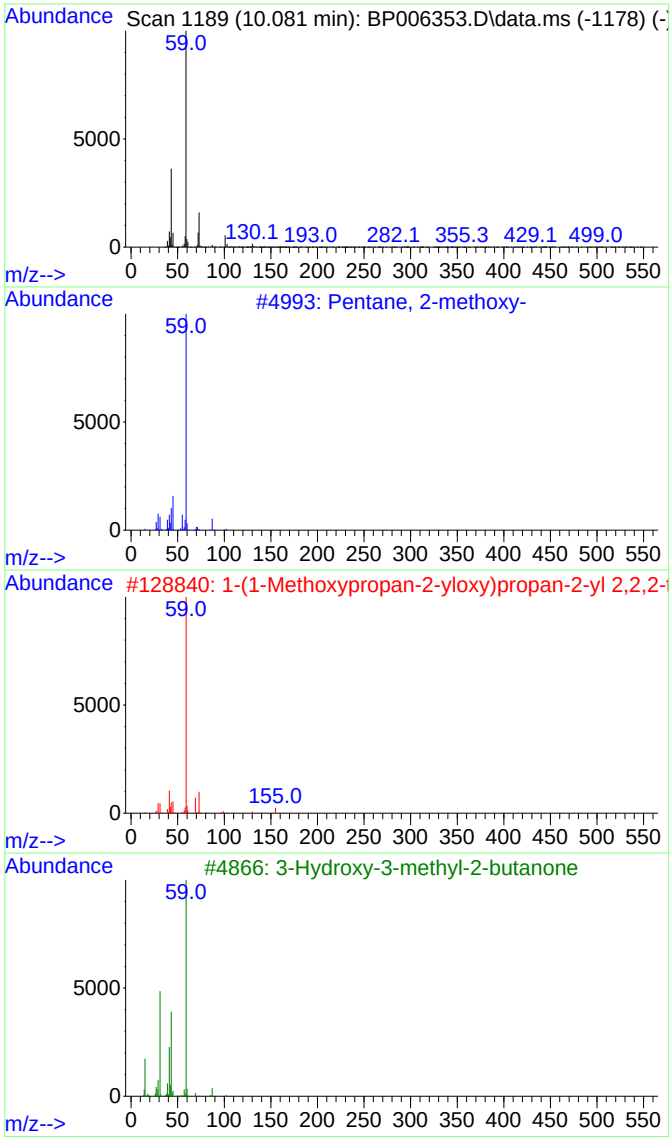
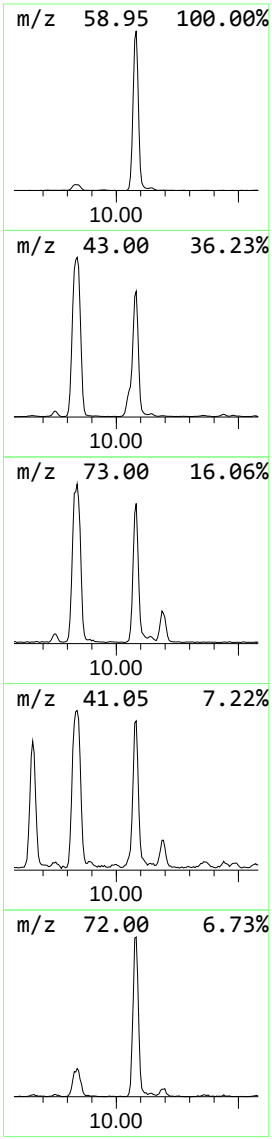
Peak Number 7 Pentane, 2-methoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.080	15.19 ng/ul	1705460	Naphthalene-d8	10.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methoxy-	102	C6H14O	006795-88-6	59
2	1-(1-Methoxypropan-2-yloxy)propa...	244	C9H15F3O4	1010367-08-7	56
3	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	53
4	2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	50
5	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	47





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Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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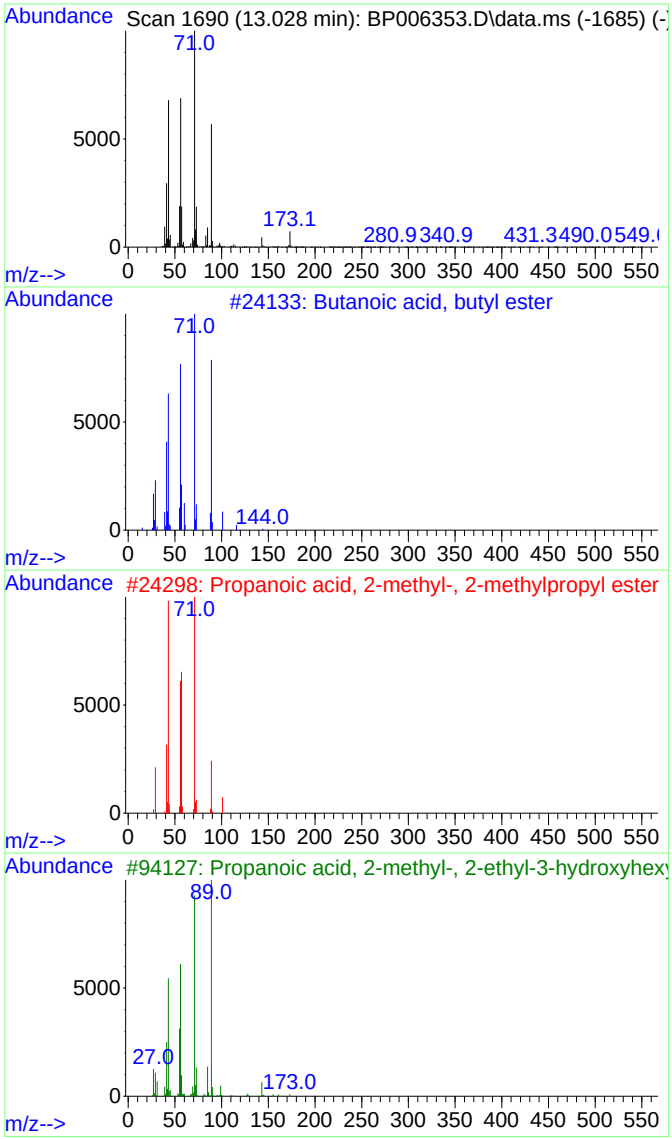
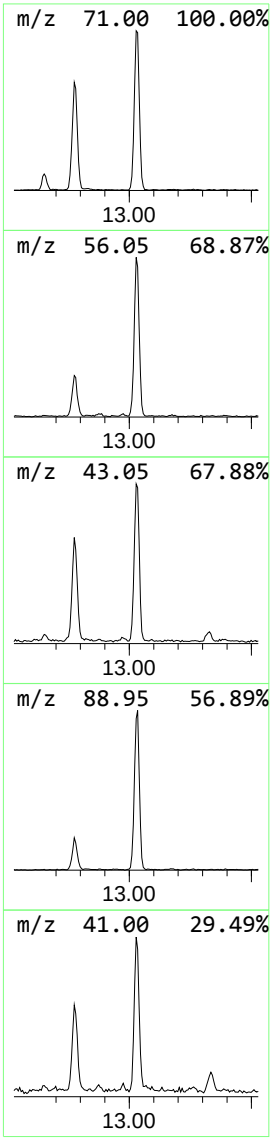
Peak Number 8 Butanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.030	2.26 ng/ul	318645	Acenaphthene-d10	14.410

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	87
2	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	64
3	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	64
4	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	64
5	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	56





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Sample : M3123-01
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ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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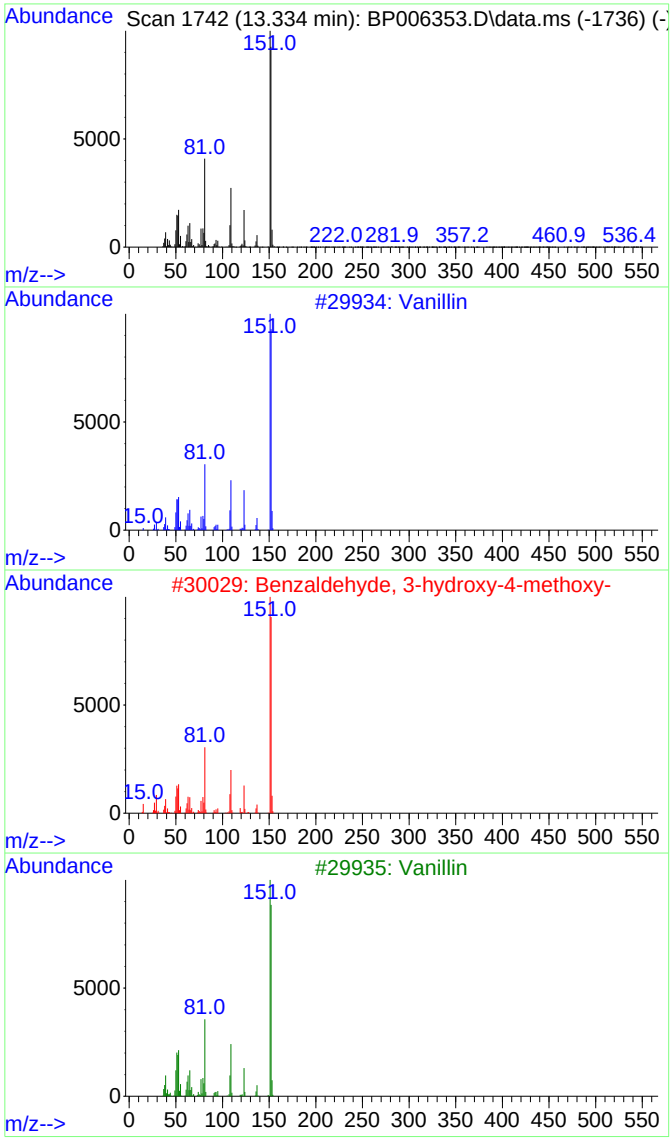
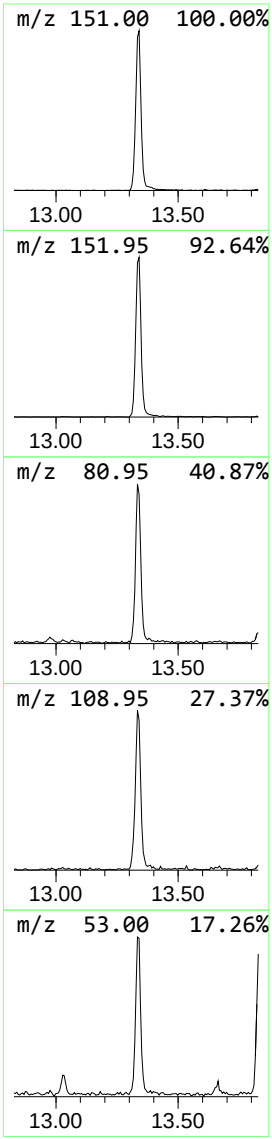
Peak Number 9 Vanillin Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.330	2.47 ng/ul	349078	Acenaphthene-d10	14.410

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	97
2	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97
3	Vanillin	152	C8H8O3	000121-33-5	96
4	Vanillin	152	C8H8O3	000121-33-5	96
5	Vanillin	152	C8H8O3	000121-33-5	95





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Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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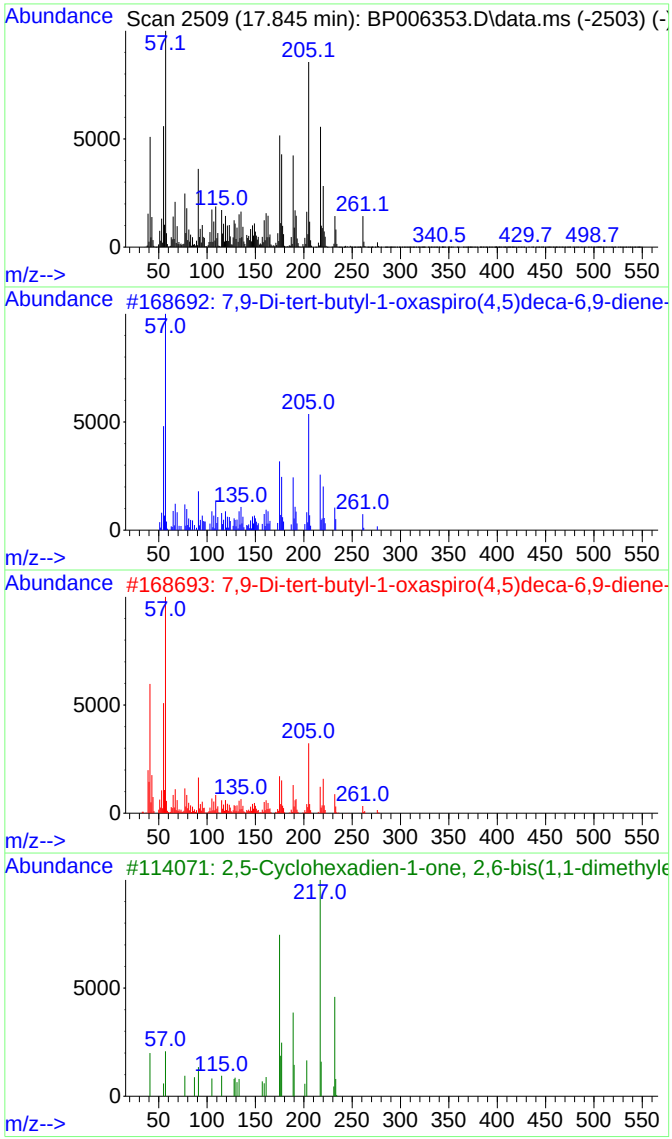
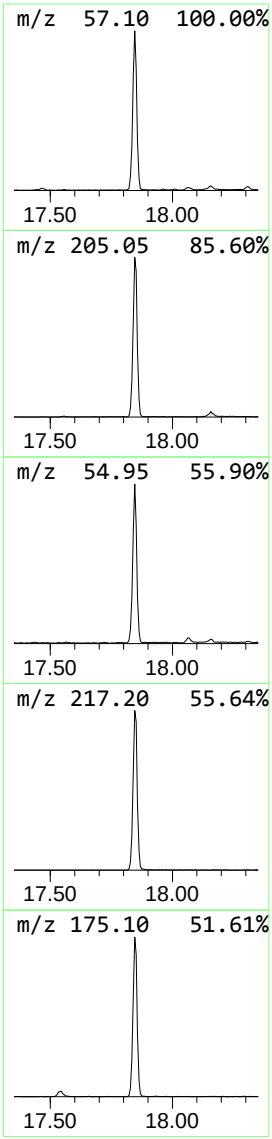
Peak Number 10 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	14.46 ng/ul	2250900	Phenanthrene-d10	17.163

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	97
3	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	80
4	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	68
5	1,3-Pentadiene, 1,1-diphenyl-, (Z)-	220	C17H16	015295-31-5	51





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Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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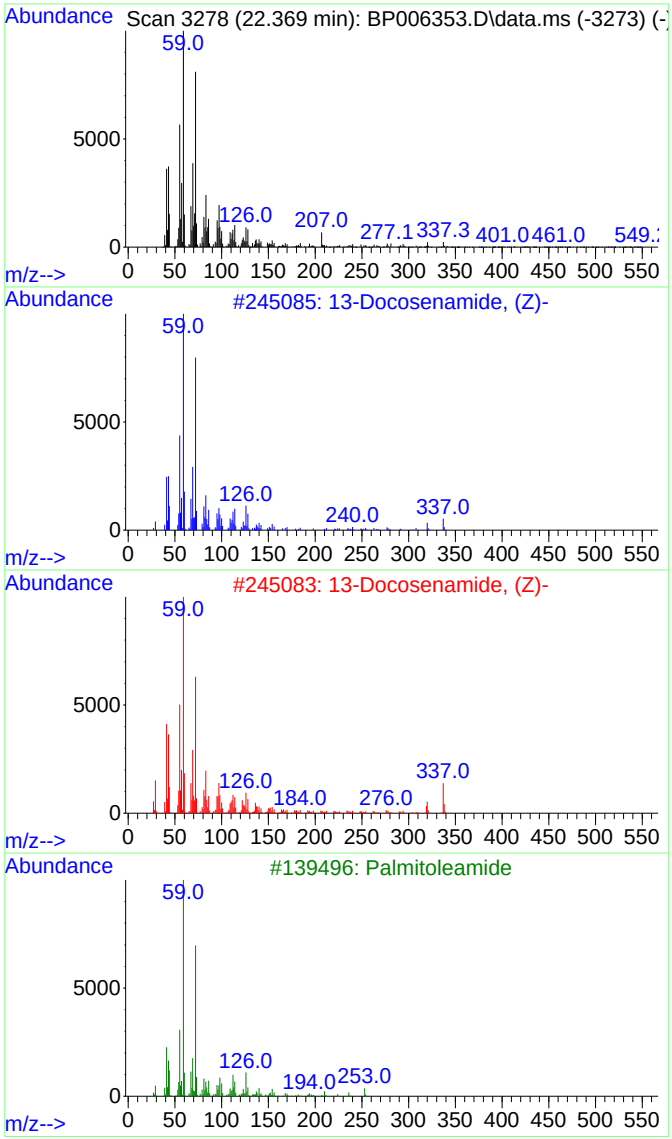
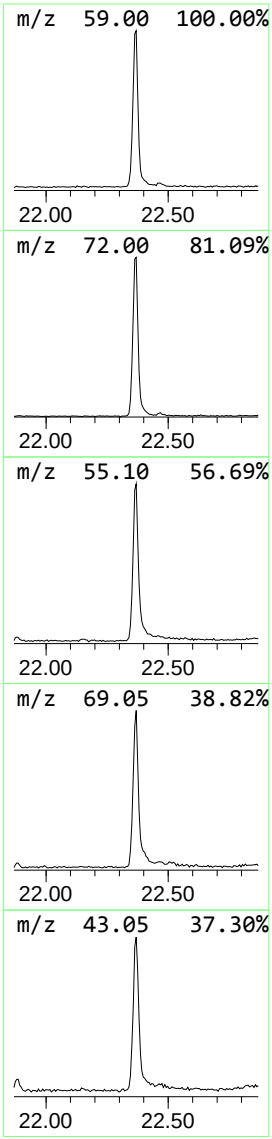
Peak Number 11 13-Docosenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.370	6.95 ng/ul	1154880	Chrysene-d12	21.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	99
2	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	74
3	Palmitoleamide	253	C16H31NO	106010-22-4	64
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	60
5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	52





Tentatively Identified Compound (LSC) summary

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Data File : BP006353.D
Acq On : 26 Jul 2021 23:44
Operator : CG/JU
Sample : M3123-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0060

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.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
Cyclohexanone	5.860	2.1	ng/ul	169630	1	7.787	1586660	20.0
Ethanol, 2-butoxy-	5.890	2.3	ng/ul	181053	1	7.787	1586660	20.0
1-Hexanol, 2-et...	7.860	3.1	ng/ul	243310	1	7.787	1586660	20.0
Phenol, 2-methoxy-	8.880	6.0	ng/ul	478894	1	7.787	1586660	20.0
Ethanol, 2-(hex...	9.120	2.3	ng/ul	179093	1	7.787	1586660	20.0
Dipropylene gly...	9.840	11.3	ng/ul	1265190	2	10.569	2245030	20.0
Pentane, 2-meth...	10.080	15.2	ng/ul	1705460	2	10.569	2245030	20.0
Butanoic acid, ...	13.030	2.3	ng/ul	318645	3	14.410	2822510	20.0
Vanillin	13.330	2.5	ng/ul	349078	3	14.410	2822510	20.0
7,9-Di-tert-but...	17.850	14.5	ng/ul	2250900	4	17.163	3113350	20.0
13-Docosenamide...	22.370	7.0	ng/ul	1154880	5	21.257	3324990	20.0