LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : A4D64

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_G\Methods\SFAM-EPA-BG070224.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG061981.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.480	29	32	38	rVB4	10367	15909	0.66%	0.063%
2	3.903	102	104	115	rBV	48760	74930	3.10%	0.297%
3	4.249	160	163	166	rBV3	3759	4557	0.19%	0.018%
4	4.608	221	224	229	rBV5	5211	6653	0.28%	0.026%
5	5.401	354	359	362	rBV5	4501	7360	0.30%	0.029%
6	5.912	445	446	452	rVB3	3849	4658	0.19%	0.018%
7	6.147	477	486	491	rBV3	39558	81453	3.37%	0.322%
8	6.711	576	582	589	rBV4	16185	29673	1.23%	0.117%
9	7.099	642	648	651	rBV3	3743	7628	0.32%	0.030%
10	7.122	651	652	656	rVB3	5257	4610	0.19%	0.018%
11	7.169	656	660	663	rBV3	4453	8068	0.33%	0.032%
12	7.240	666	672	680	rVV	90305	170693	7.06%	0.675%
13	7.416	694	702	711	rBV	206489	350864	14.52%	1.388%
14	7.622	731	737	745	rBV	263157	449093	18.58%	1.777%
15	7.839	768	774	775	rBV3	7413	10375	0.43%	0.041%
16	8.092	810	817	822	rBV2	171245	301889	12.49%	1.195%
17	8.139	822	825	830	rVB	41192	60489	2.50%	0.239%
18	8.233	838	841	845	rBV5	5847	7866	0.33%	0.031%
19	8.327	852	857	861	rBV3	10011	16174	0.67%	0.064%
20	8.603	900	904	907	rBV2	3788	6537	0.27%	0.026%
21	8.685	910	918	922	rBV4	9093	18843	0.78%	0.075%
22	8.791	929	936	944	rBV	253674	440725	18.23%	1.744%
23	8.867	944	949	954	rVV2	45654	67652	2.80%	0.268%
24	8.920	954	958	962	rVV3	12484	19029	0.79%	0.075%
25	9.185	998	1003	1009	rBV4	29866	52993	2.19%	0.210%
26	9.261	1009	1016	1023	rVV	338741	592088	24.50%	2.343%
27	9.331	1023	1028	1034	rVB3	70043	108405	4.49%	0.429%
28	9.408	1034	1041	1047	rBV2	37443	68633	2.84%	0.272%
29	9.637	1076	1080	1084	rBV2	6698	9924	0.41%	0.039%
30	9.984	1129	1139	1149	rBV	309437	570102	23.59%	2.256%
31	10.119	1156	1162	1172	rBV4	33427	96249	3.98%	0.381%
32	10.219	1174	1179	1185	rVB5	11844	24988	1.03%	0.099%
33	10.360	1197	1203	1210	rVB2	174080	280768	11.62%	1.111%
34	10.518	1223	1230	1238	rBV	436451	800166	33.11%	3.166%
35	10.671	1246	1256	1264	rBV5	16279	43867	1.81%	0.174%
36	10.730	1264	1266	1267	rBV2	7914	6416	0.27%	0.025%

LSC Area Percent Report

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Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

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_G\Methods\SFAM-EPA-BG070224.MA.M
Title : SVOA CALIBRATION

37	10.800	1274	1278	1285	rVB6	7555	14182	0.59%	0.056%
38	10.906	1289	1296	1303	rBV	295122	533694	22.08%	2.112%
39	11.035	1311	1318	1325	rBV	235400	419072	17.34%	1.658%
40	11.176	1339	1342	1352	rVB2	16946	30081	1.24%	0.119%
41	11.417	1379	1383	1389	rVB3	9940	13451	0.56%	0.053%
42	11.482	1390	1394	1398	rVV5	8779	15969	0.66%	0.063%
43	11.558	1402	1407	1413	rVB4	12530	23661	0.98%	0.094%
44	11.688	1423	1429	1435	rBV4	27384	49999	2.07%	0.198%
45	12.058	1486	1492	1495	rBV5	7313	14779	0.61%	0.058%
46	12.122	1500	1503	1510	rVB4	7803	10823	0.45%	0.043%
47	12.304	1529	1534	1539	rVB5	7903	14066	0.58%	0.056%
48	12.481	1558	1564	1568	rVB2	8890	15911	0.66%	0.063%
49	12.610	1585	1586	1591	rVB3	5752	5788	0.24%	0.023%
50	12.839	1621	1625	1627	rVB3	5367	5650	0.23%	0.022%
51	13.074	1659	1665	1672	rBV2	31738	60279	2.49%	0.239%
52	13.262	1691	1697	1701	rBV2	12717	20499	0.85%	0.081%
53	13.315	1701	1706	1713	rBV	48266	74051	3.06%	0.293%
54	13.515	1736	1740	1741	rBV2	3735	5458	0.23%	0.022%
55	13.550	1744	1746	1749	rVV2	5650	7140	0.30%	0.028%
56	13.597	1749	1754	1757	rVV	127547	217233	8.99%	0.860%
57	13.626	1757	1759	1767	rVB3	135914	201419	8.33%	0.797%
58	13.944	1809	1813	1822	rVB6	25496	50047	2.07%	0.198%
59	14.114	1836	1842	1849	rBV	900707	1316301	54.46%	5.209%
60	14.238	1858	1863	1868	rBV4	20194	40310	1.67%	0.160%
61	14.414	1886	1893	1900	rVB	880180	1415401	58.56%	5.601%
62	14.473	1900	1903	1908	rVB3	4015	6670	0.28%	0.026%
63	14.520	1908	1911	1914	rBV3	6432	8809	0.36%	0.035%
64	14.567	1915	1919	1921	rBV4	4588	6272	0.26%	0.025%
65	14.637	1926	1931	1936	rBV2	38435	62962	2.61%	0.249%
66	14.713	1937	1944	1951	rVB	513384	825778	34.17%	3.268%
67	14.784	1951	1956	1959	rBV3	6921	9441	0.39%	0.037%
68	14.907	1968	1977	1986	rBV2	111491	207456	8.58%	0.821%
69	15.125	2009	2014	2015	rBV4	9757	12648	0.52%	0.050%
70	15.201	2024	2027	2029	rBV2	3694	5110	0.21%	0.020%
71	15.389	2056	2059	2063	rVV4	6673	10948	0.45%	0.043%
72	15.442	2065	2068	2073	rVB4	12286	12384	0.51%	0.049%
73	15.706	2105	2113	2120	rBV	1178128	1796652	74.34%	7.110%
74	15.818	2125	2132	2138	rBV	514258	765286	31.66%	3.028%
75	15.941	2149	2153	2158	rVB5	8885	13862	0.57%	0.055%
76	16.059	2170	2173	2177	rBV5	8242	12045	0.50%	0.048%

LSC Area Percent Report

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Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

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

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Title : SVOA CALIBRATION

77	16.447	2236	2239	2243	rBV2	15623	19661	0.81%	0.078%
78	16.934	2320	2322	2327	rVB5	9760	11751	0.49%	0.047%
79	17.034	2337	2339	2345	rVB4	8674	10768	0.45%	0.043%
80	17.134	2353	2356	2359	rVB	3656	4535	0.19%	0.018%
81	17.457	2404	2411	2417	rVB	641134	974107	40.30%	3.855%
82	17.557	2422	2428	2435	rVB	1345984	2037577	84.30%	8.063%
83	17.722	2451	2456	2458	rBV5	5276	8642	0.36%	0.034%
84	17.822	2469	2473	2476	rBV3	4833	8527	0.35%	0.034%
85	18.109	2516	2522	2528	rBV	1269391	1744149	72.16%	6.902%
86	18.585	2599	2603	2607	rBV4	10922	16735	0.69%	0.066%
87	18.673	2613	2618	2619	rBV5	7824	9022	0.37%	0.036%
88	18.820	2641	2643	2646	rBV3	4996	5786	0.24%	0.023%
89	19.402	2735	2742	2747	rBV3	19307	31202	1.29%	0.123%
90	19.531	2761	2764	2767	rVB3	4178	4814	0.20%	0.019%
91	19.631	2779	2781	2784	rVB4	5083	5673	0.23%	0.022%
92	19.731	2793	2798	2802	rBV4	26093	42805	1.77%	0.169%
93	19.837	2809	2816	2821	rBV	1519875	2248357	93.03%	8.897%
94	20.806	2979	2981	2984	rBV4	8745	9171	0.38%	0.036%
95	21.717	3129	3136	3142	rVB	570971	979458	40.52%	3.876%
96	23.180	3380	3385	3394	rVB4	94551	212949	8.81%	0.843%
97	24.596	3624	3626	3631	rBV5	9645	13273	0.55%	0.053%
98	24.731	3638	3649	3663	rVB2	866220	2416934	100.00%	9.564%
99	24.954	3677	3687	3701	rVB	416513	1213519	50.21%	4.802%
100	26.106	3877	3883	3892	rVB8	31751	88608	3.67%	0.351%

Sum of corrected areas: 25269937

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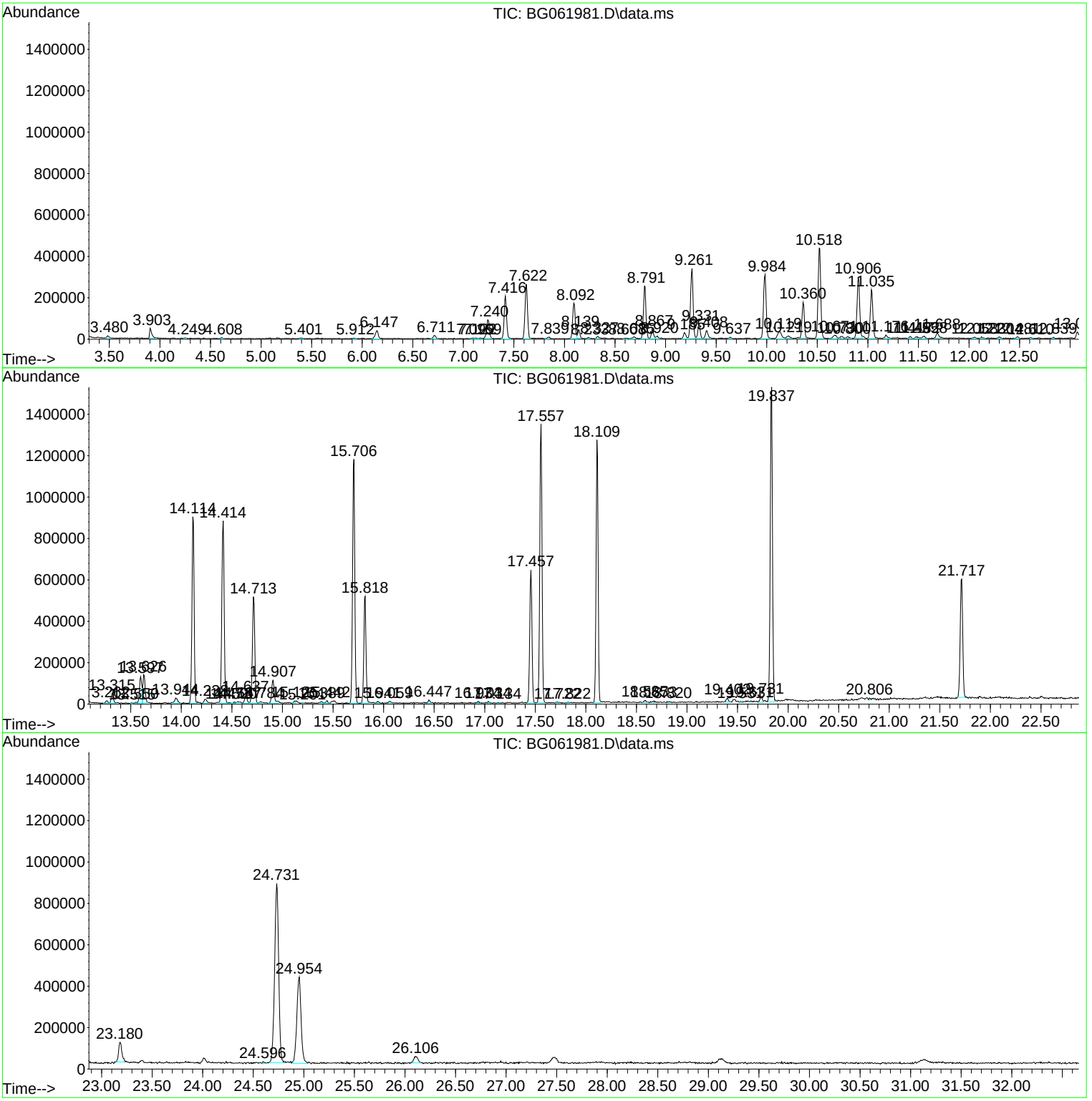
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
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Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

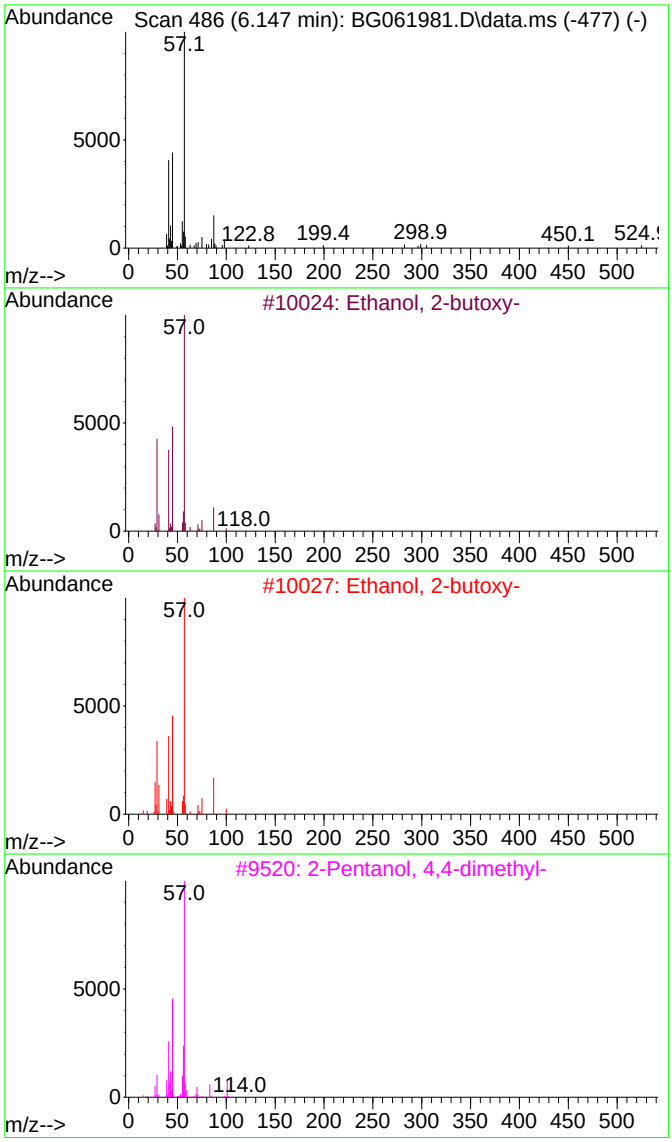
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

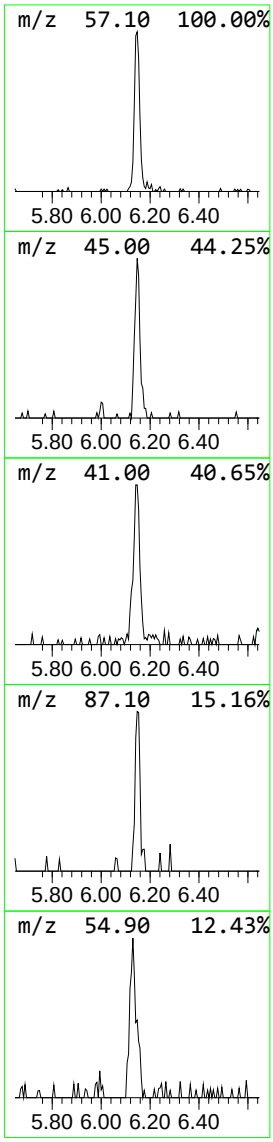
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.147	5.40 ng/ul	81453	1,4-Dichlorobenzene-d4	8.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	59
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	40
3	2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	33
4	2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	33
5	3-Pentanol, 3-(1,1-dimethylethyl...	200	C13H28O	041902-42-5	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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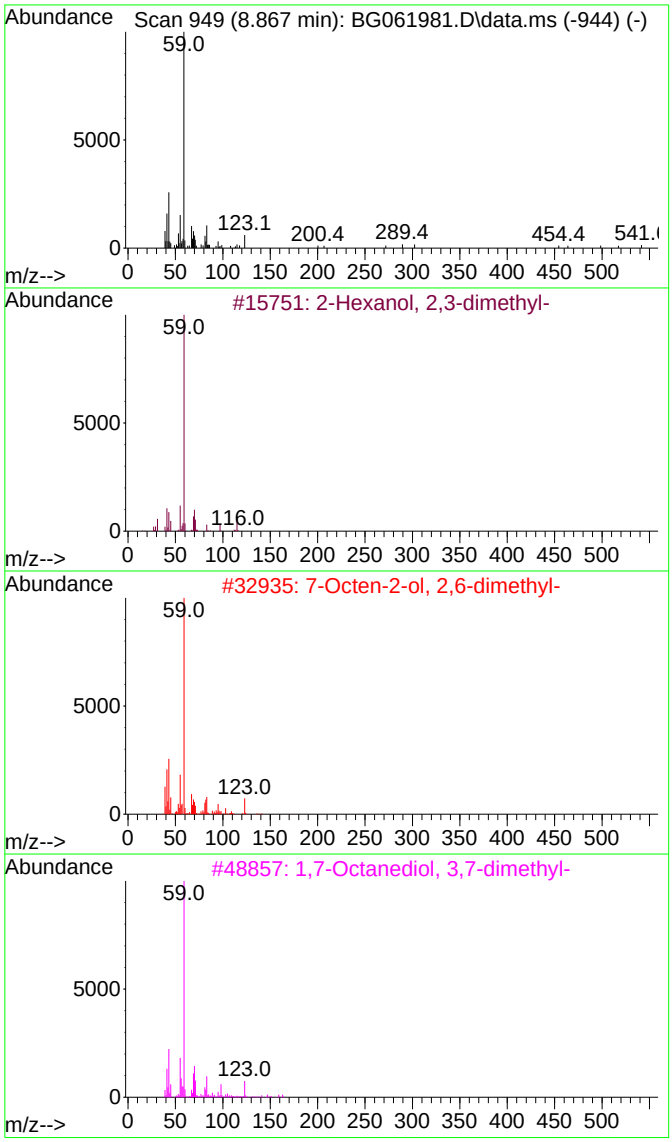
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

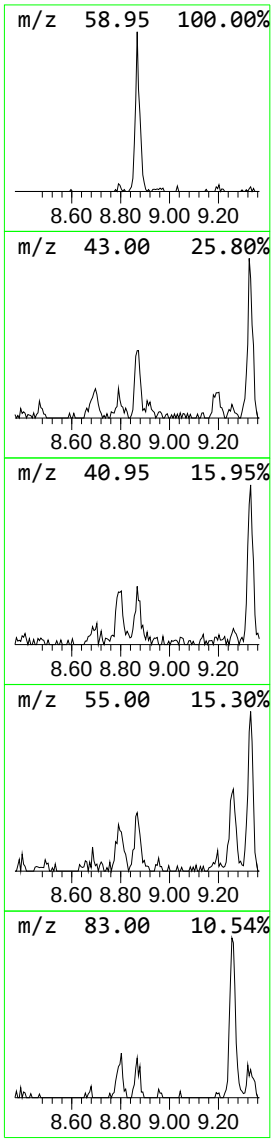
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Hexanol, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.867	4.48 ng/ul	67652	1,4-Dichlorobenzene-d4	8.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	59
2	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	50
3	1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4	50
4	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	50
5	2-Hexanol, 2,5-dimethyl-, (S)-	130	C8H18O	003730-60-7	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

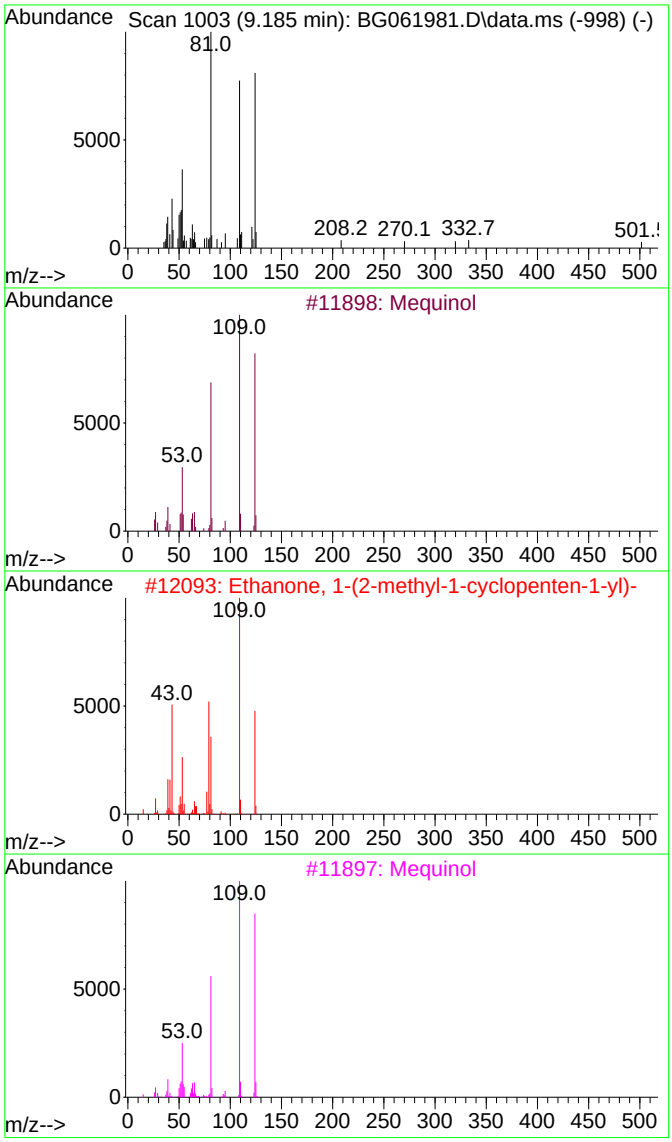
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

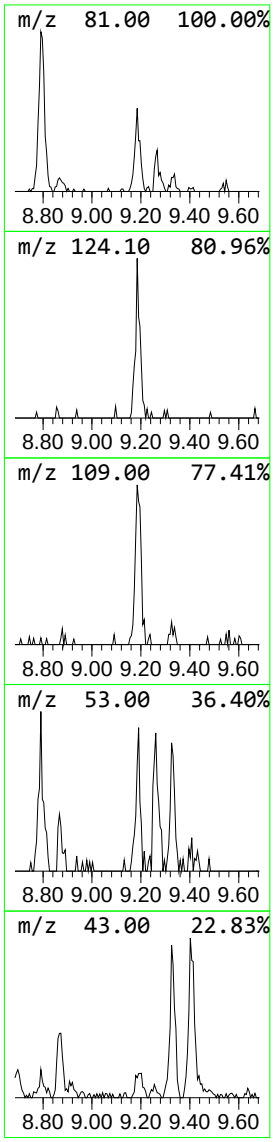
****

Peak Number 3 Mequinol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.185	3.51 ng/ul	52993	1,4-Dichlorobenzene-d4	8.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mequinol	124	C7H8O2	000150-76-5	64
2	Ethanone, 1-(2-methyl-1-cyclopen...	124	C8H12O	003168-90-9	60
3	Mequinol	124	C7H8O2	000150-76-5	52
4	Mequinol	124	C7H8O2	000150-76-5	52
5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	49





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Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

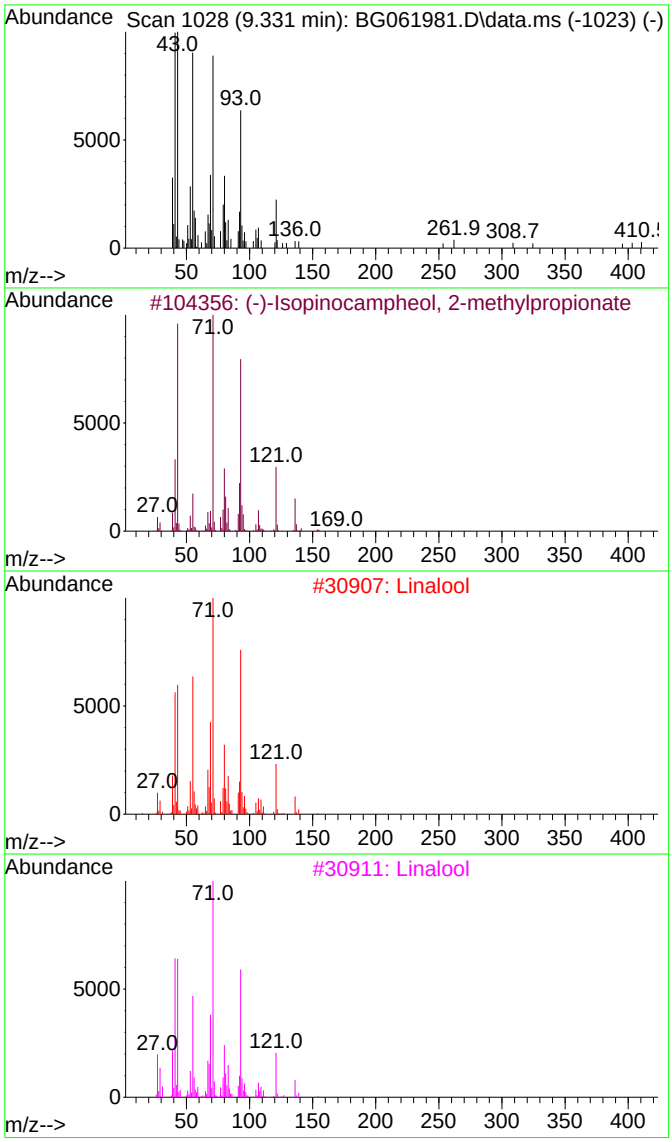
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

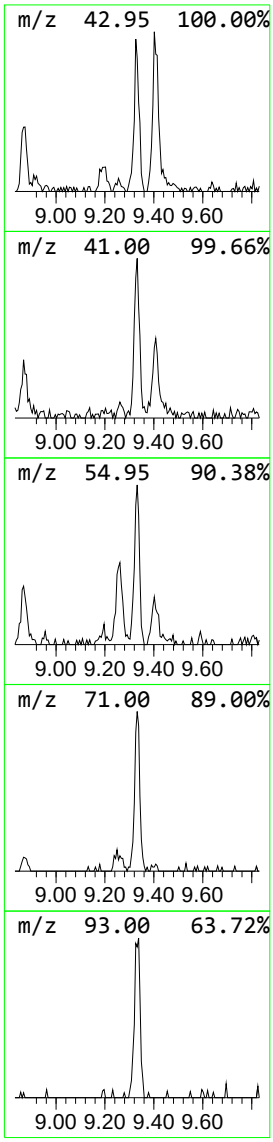
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (-)-Isopinocampheol, 2-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.331	7.18 ng/ul	108405	1,4-Dichlorobenzene-d4	8.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(-)-Isopinocampheol, 2-methylpro...	224	C14H24O2	1000462-98-0	86
2	Linalool	154	C10H18O	000078-70-6	78
3	Linalool	154	C10H18O	000078-70-6	64
4	1,5-Dimethyl-1-vinyl-4-hexenyl b...	224	C14H24O2	000078-36-4	47
5	1,6-Octadien-3-ol, 3,7-dimethyl-...	182	C11H18O2	000115-99-1	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

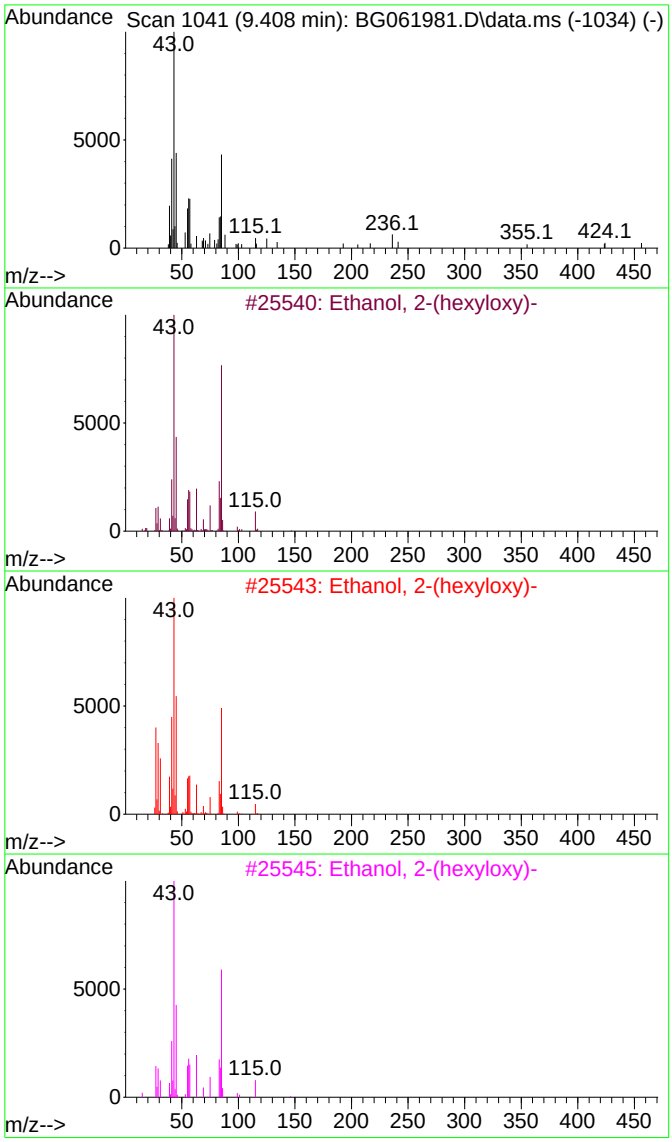
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

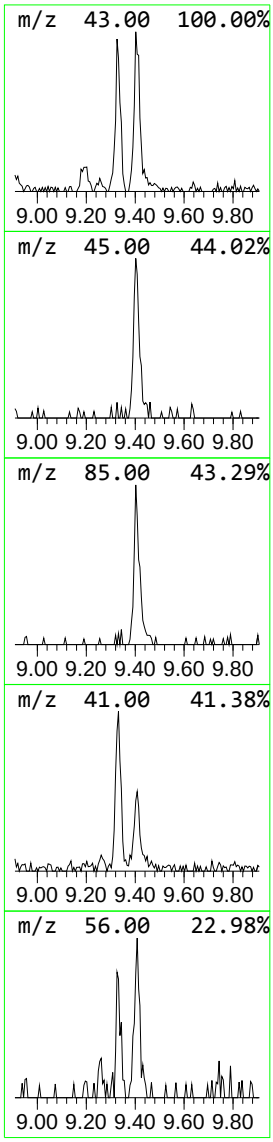
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-(hexyloxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.408	4.55 ng/ul	68633	1,4-Dichlorobenzene-d4	8.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	72
2	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	72
3	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	56
4	Oxalic acid, isohexyl propyl ester	216	C11H20O4	1000309-32-6	43
5	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

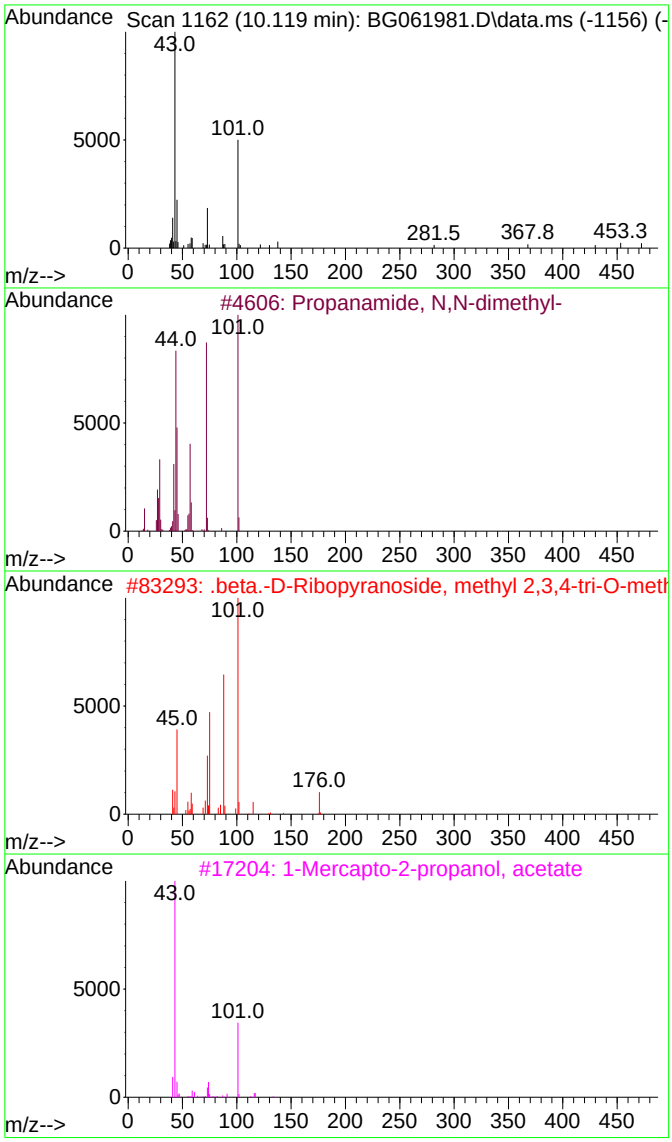
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

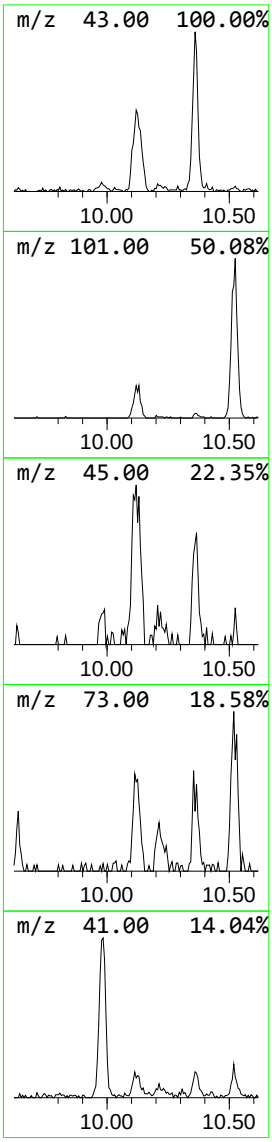
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.119	3.61 ng/ul	96249	Naphthalene-d8	10.906

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanamide, N,N-dimethyl-	101	C5H11NO	000758-96-3	40
2	.beta.-D-Ribopyranoside, methyl ...	206	C9H18O5	002876-87-1	39
3	1-Mercapto-2-propanol, acetate	134	C5H10O2S	035508-06-6	28
4	3H-1,2,4-Triazole-3-thione, 1,2-...	101	C2H3N3S	003179-31-5	25
5	Glutaric acid, heptyl 4-methoxy-...	330	C18H34O5	1000359-41-6	25





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Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

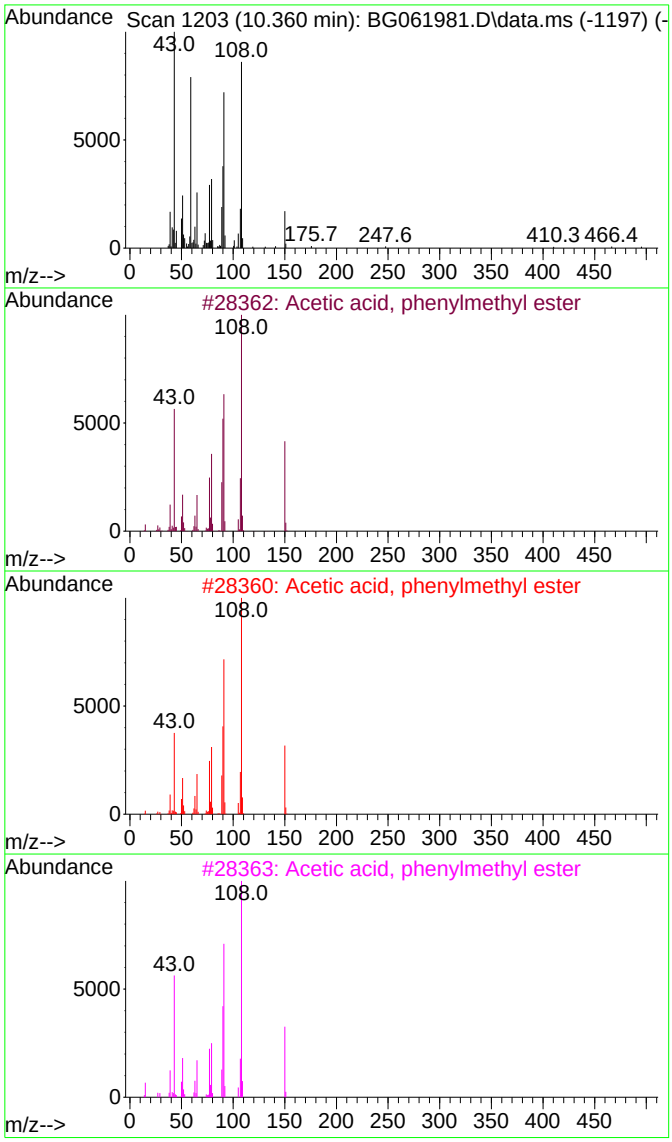
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

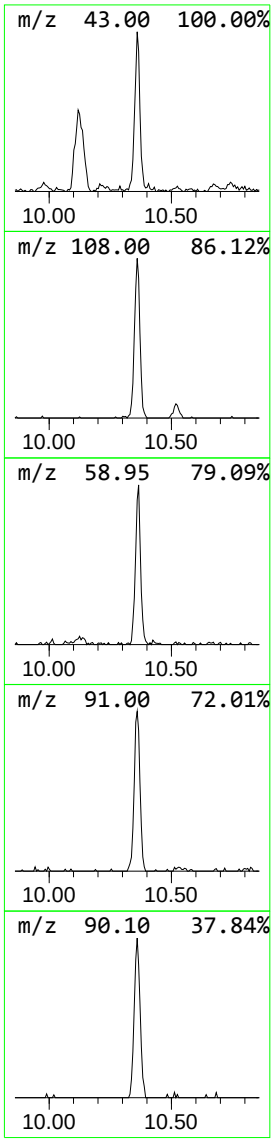
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Acetic acid, phenylmethyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	10.52 ng/ul	280768	Naphthalene-d8	10.906

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	92
2	Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	83
3	Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	50
4	Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	43
5	Phenol, 2-methyl-	108	C7H8O	000095-48-7	43





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Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

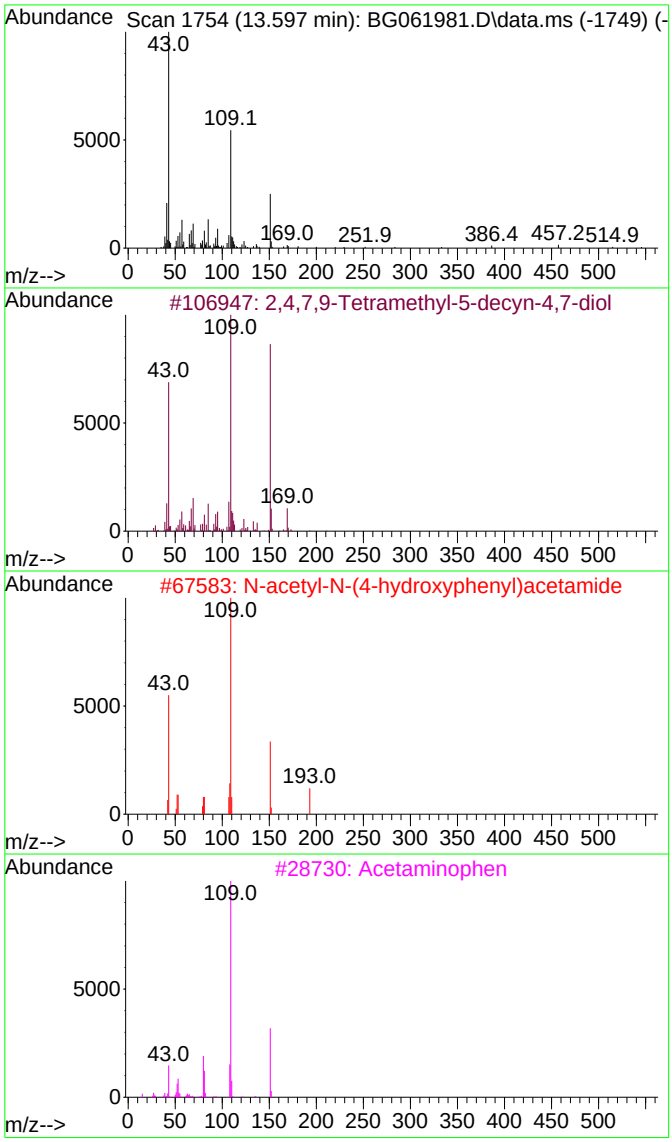
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

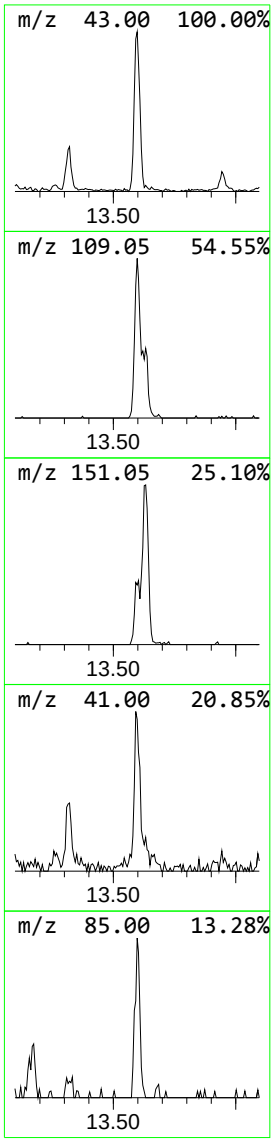
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.597	5.26 ng/ul	217233	Acenaphthene-d10	14.719

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	78
2	N-acetyl-N-(4-hydroxyphenyl)acet...	193	C10H11NO3	1000389-51-9	53
3	Acetaminophen	151	C8H9NO2	000103-90-2	52
4	Metacetamol	151	C8H9NO2	000621-42-1	52
5	Metacetamol	151	C8H9NO2	000621-42-1	52





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Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

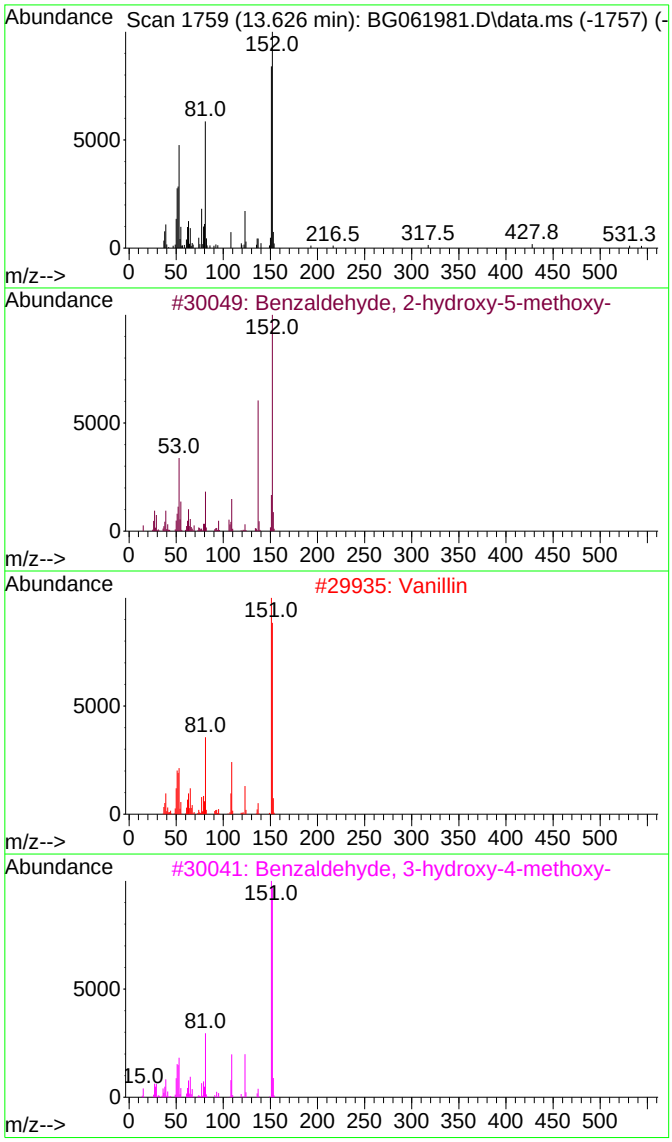
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

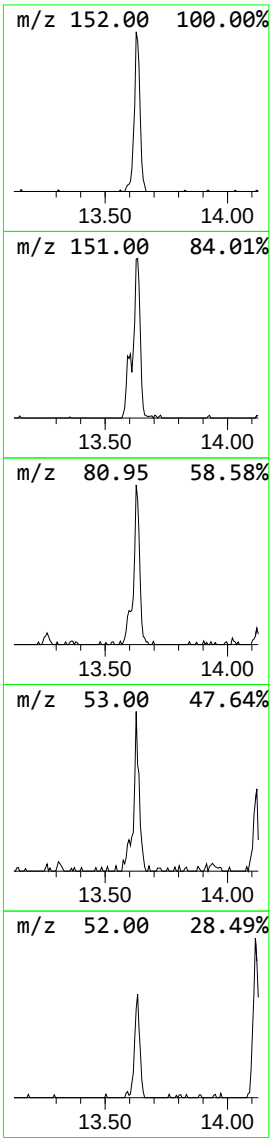
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzaldehyde, 2-hydroxy-5-m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.627	4.88 ng/ul	201419	Acenaphthene-d10	14.719

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 2-hydroxy-5-methoxy-	152	C8H8O3	000672-13-9	50
2	Vanillin	152	C8H8O3	000121-33-5	50
3	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	38
4	Vanillin	152	C8H8O3	000121-33-5	38
5	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	38





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Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

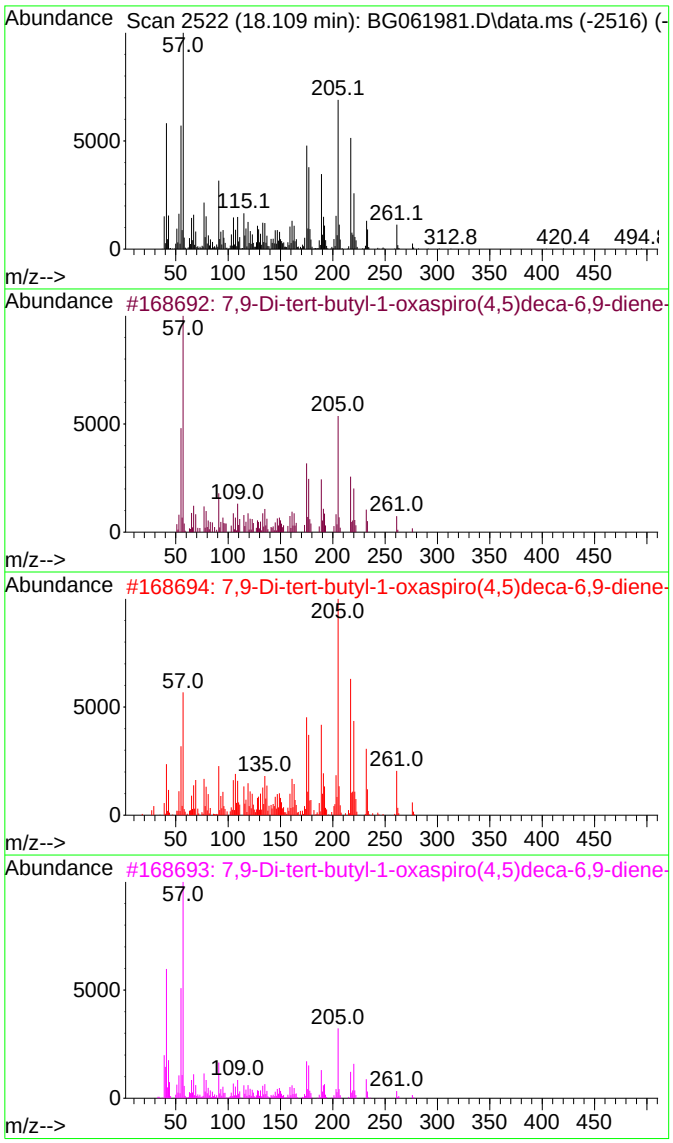
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

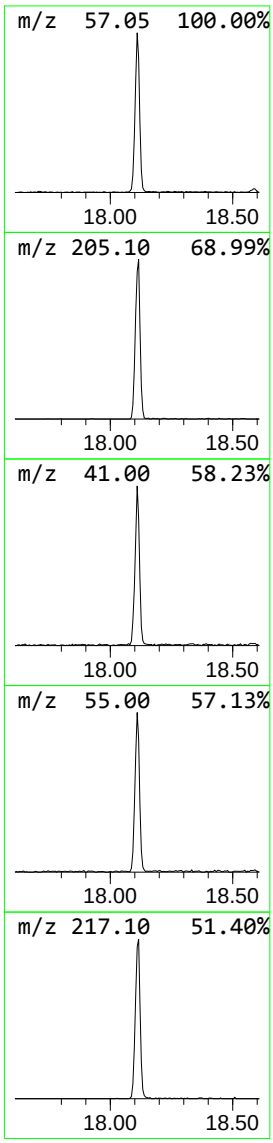
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.109	35.81 ng/ul	1744150	Phenanthrene-d10	17.457

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	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	86
4	2,4-Diamino-6-nitroquinazoline	205	C8H7N5O2	007154-34-9	42
5	(Z)-((2-Benzylideneheptyl)oxy)tr...	276	C17H28OSi	1000495-65-9	41





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Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

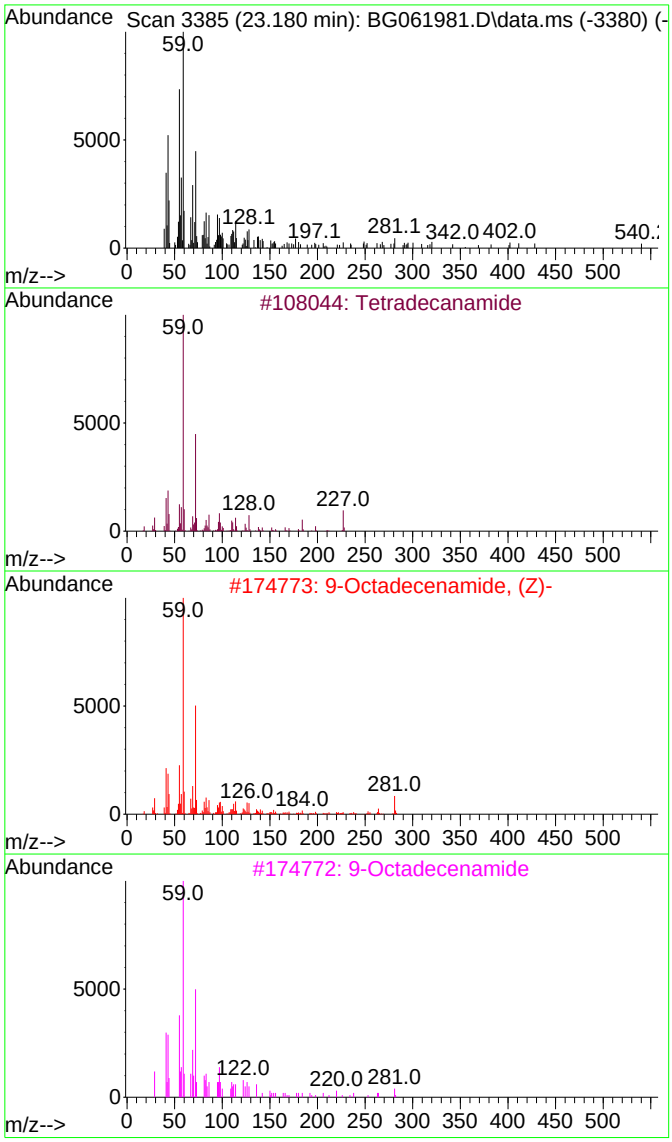
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

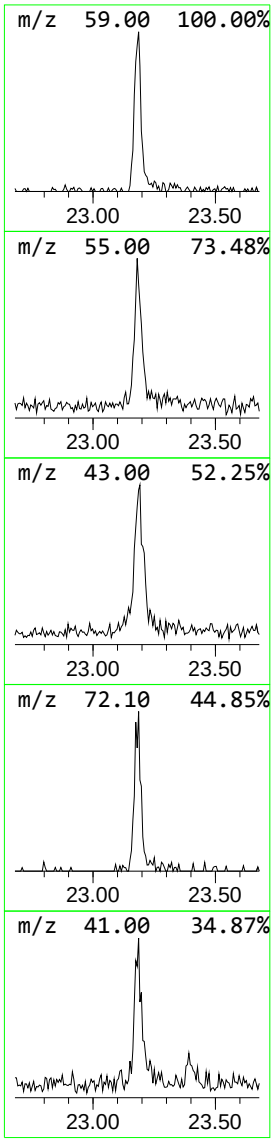
****

Peak Number 11 Tetradecanamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.180	4.35 ng/ul	212949	Chrysene-d12	21.717

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanamide	227	C14H29NO	000638-58-4	94
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
3	9-Octadecenamide	281	C18H35NO	003322-62-1	87
4	Tetradecanamide	227	C14H29NO	000638-58-4	74
5	Hexadecanamide	255	C16H33NO	000629-54-9	72





Tentatively Identified Compound (LSC) summary

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Data File : BG061981.D
Acq On : 9 Jul 2024 20:03
Operator : MA/JU
Sample : P3091-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4D64

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.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
Ethanol, 2-butoxy-	6.147	5.4	ng/ul	81453	1	8.098	301889	20.0
2-Hexanol, 2,3-...	8.867	4.5	ng/ul	67652	1	8.098	301889	20.0
Mequinol	9.185	3.5	ng/ul	52993	1	8.098	301889	20.0
(-)-Isopinocamp...	9.331	7.2	ng/ul	108405	1	8.098	301889	20.0
Ethanol, 2-(hex...	9.408	4.5	ng/ul	68633	1	8.098	301889	20.0
unknown-01	10.119	3.6	ng/ul	96249	2	10.906	533694	20.0
Acetic acid, ph...	10.360	10.5	ng/ul	280768	2	10.906	533694	20.0
2,4,7,9-Tetrame...	13.597	5.3	ng/ul	217233	3	14.719	825778	20.0
Benzaldehyde, 2...	13.627	4.9	ng/ul	201419	3	14.719	825778	20.0
7,9-Di-tert-but...	18.109	35.8	ng/ul	1744150	4	17.457	974107	20.0
Tetradecanamide	23.180	4.3	ng/ul	212949	5	21.717	979458	20.0