LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

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\SOM-EPA-BP022720MA.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	2.946	7	10	21	rVB	120568	148252	1.01%	0.056%
2	3.146	40	44	51	rBV	84936	111952	0.76%	0.042%
3	3.211	51	55	63	rVB	65193	84781	0.58%	0.032%
4	3.493	100	103	110	rVB2	22278	28430	0.19%	0.011%
5	3.875	164	168	171	rBV	28855	41388	0.28%	0.016%
6	4.140	207	213	218	rBV	31220	61090	0.42%	0.023%
7	4.193	219	222	233	rVB	26729	51874	0.35%	0.020%
8	4.393	252	256	264	rVB	7108	16420	0.11%	0.006%
9	4.775	316	321	332	rBV	177882	252266	1.72%	0.095%
10	4.975	350	355	365	rVB	173489	263629	1.79%	0.099%
11	5.840	497	502	511	rVB	50930	78941	0.54%	0.030%
12	6.528	612	619	627	rBV2	8990	18245	0.12%	0.007%
13	6.775	654	661	662	rBV	84705	122805	0.84%	0.046%
14	6.811	662	667	681	rVB	1397199	2205281	15.00%	0.830%
15	6.964	687	693	704	rBV	1128255	1772286	12.05%	0.667%
16	7.164	719	727	729	rBV	1554964	2340545	15.92%	0.881%
17	7.193	729	732	744	rVV	1753780	2674751	18.19%	1.006%
18	7.622	797	805	819	rBV	994291	1602467	10.90%	0.603%
19	8.158	887	896	902	rBV	1423545	3388178	23.04%	1.275%
20	8.211	902	905	914	rVB	158980	235766	1.60%	0.089%
21	8.334	919	926	936	rVV	1731992	2727310	18.55%	1.026%
22	8.434	936	943	965	rVB	2038463	3233726	21.99%	1.217%
23	8.693	980	987	993	rBV	822265	1331681	9.06%	0.501%
24	8.769	993	1000	1003	rVV	1233980	2061578	14.02%	0.776%
25	8.811	1003	1007	1020	rVB	2199329	3427691	23.31%	1.290%
26	9.093	1047	1055	1063	rBV	2867050	4497250	30.58%	1.692%
27	9.487	1114	1122	1124	rBV	1026455	1635850	11.12%	0.616%
28	9.516	1124	1127	1134	rVV	1481276	2329393	15.84%	0.876%
29	9.587	1134	1139	1151	rVB2	86729	232572	1.58%	0.088%
30	10.028	1206	1214	1238	rBV2	1773894	4937571	33.58%	1.858%
31	10.246	1247	1251	1256	rBV2	10657	18931	0.13%	0.007%
32	10.310	1256	1262	1269	rVV	78302	150378	1.02%	0.057%
33	10.399	1269	1277	1281	rVV	1344667	2341840	15.93%	0.881%
34	10.446	1281	1285	1293	rVV	1207632	1955083	13.30%	0.736%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

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\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

35	10.534	1293	1300	1314	rVV	362707	693164	4.71%	0.261%
36	11.122	1394	1400	1406	rBV	81078	135661	0.92%	0.051%
37	11.693	1488	1497	1519	rBV	2476215	4084568	27.78%	1.537%
38	11.940	1531	1539	1547	rBV	2573836	4301522	29.25%	1.618%
39	12.193	1576	1582	1586	rBV2	28051	50664	0.34%	0.019%
40	12.287	1590	1598	1613	rVB	3979427	6454134	43.89%	2.428%
41	12.446	1615	1625	1633	rBV	1450263	2295717	15.61%	0.864%
42	12.687	1659	1666	1672	rBV	1961049	2883017	19.61%	1.085%
43	12.734	1672	1674	1686	rVB	58324	107202	0.73%	0.040%
44	13.093	1721	1735	1747	rBV	4339637	6261442	42.58%	2.356%
45	13.334	1763	1776	1792	rBV	2227376	3344297	22.74%	1.258%
46	13.469	1792	1799	1808	rVB	99654	149234	1.01%	0.056%
47	13.681	1824	1835	1842	rBV	2807005	4035243	27.44%	1.518%
48	13.757	1842	1848	1855	rVV	926309	1314338	8.94%	0.495%
49	13.840	1855	1862	1874	rVB	2040213	2851474	19.39%	1.073%
50	13.951	1874	1881	1883	rBV	3566967	5437175	36.98%	2.046%
51	13.981	1883	1886	1895	rVB	5096600	7160534	48.70%	2.694%
52	14.263	1927	1934	1940	rBV	2097884	3079656	20.94%	1.159%
53	14.328	1940	1945	1958	rVB	523582	771159	5.24%	0.290%
54	14.487	1963	1972	1990	rBV2	2754013	5337607	36.30%	2.008%
55	14.634	1990	1997	1999	rVV	2482363	3507729	23.86%	1.320%
56	14.663	1999	2002	2009	rVB	2032321	2897774	19.71%	1.090%
57	14.910	2036	2044	2054	rBV	608872	968345	6.59%	0.364%
58	15.104	2070	2077	2089	rBV	4490582	5817927	39.57%	2.189%
59	15.263	2097	2104	2109	rBV	4553102	6523350	44.36%	2.454%
60	15.316	2109	2113	2119	rVV	1045446	1435487	9.76%	0.540%
61	15.387	2119	2125	2138	rVV	1063877	1485371	10.10%	0.559%
62	15.534	2138	2150	2165	rVB	5848914	8642900	58.78%	3.252%
63	15.769	2184	2190	2209	rBV	223294	480234	3.27%	0.181%
64	15.910	2209	2214	2220	rVB	69830	98571	0.67%	0.037%
65	16.045	2234	2237	2243	rVB5	19489	28117	0.19%	0.011%
66	16.151	2250	2255	2258	rBV	28039	40684	0.28%	0.015%
67	16.210	2258	2265	2269	rVV	73864	115945	0.79%	0.044%
68	16.257	2269	2273	2276	rVV	53056	74043	0.50%	0.028%
69	16.298	2276	2280	2282	rVV	1021926	1340859	9.12%	0.505%
70	16.334	2282	2286	2292	rVB	3738958	5270212	35.84%	1.983%
71	16.675	2337	2344	2362	rBV	2611167	3732949	25.39%	1.405%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

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\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

72	16.798	2362	2365	2369	rVV	19589	27333	0.19%	0.010%
73	17.016	2395	2402	2405	rBV	2595521	3713855	25.26%	1.397%
74	17.057	2405	2409	2414	rVV	4713902	6531511	44.42%	2.458%
75	17.116	2414	2419	2427	rVV	4904681	6914760	47.03%	2.602%
76	17.181	2427	2430	2440	rVB2	35163	69255	0.47%	0.026%
77	17.375	2458	2463	2470	rBV5	17953	28301	0.19%	0.011%
78	17.998	2558	2569	2580	rBV	5077982	5955410	40.50%	2.241%
79	18.269	2610	2615	2619	rBV2	44996	54868	0.37%	0.021%
80	19.010	2736	2741	2746	rVV	54408	73901	0.50%	0.028%
81	19.245	2776	2781	2793	rBV2	463244	654407	4.45%	0.246%
82	19.363	2795	2801	2804	rVV	6294160	9512960	64.70%	3.579%
83	19.386	2804	2805	2817	rVB	3188127	2988800	20.33%	1.125%
84	19.604	2837	2842	2851	rBV	751102	920916	6.26%	0.347%
85	19.751	2863	2867	2873	rVB3	39530	52690	0.36%	0.020%
86	20.375	2968	2973	2980	rBV	354135	443692	3.02%	0.167%
87	20.480	2988	2991	2995	rBV3	25372	31148	0.21%	0.012%
88	20.539	2997	3001	3007	rBV5	27716	51724	0.35%	0.019%
89	21.039	3082	3086	3090	rBV	140925	196681	1.34%	0.074%
90	21.098	3090	3096	3108	rVV2	9124410	14704273	100.00%	5.533%
91	21.375	3139	3143	3154	rBV	510665	662862	4.51%	0.249%
92	21.533	3165	3170	3187	rBV	8568310	9863507	67.08%	3.711%
93	21.657	3187	3191	3198	rBV	346267	475685	3.24%	0.179%
94	21.910	3229	3234	3253	rBV	4004543	5095051	34.65%	1.917%
95	22.657	3354	3361	3372	rBV	5518637	8848153	60.17%	3.329%
96	23.174	3442	3449	3454	rBV	4815370	9512783	64.69%	3.579%
97	23.222	3454	3457	3466	rVV	2893223	4835265	32.88%	1.819%
98	23.316	3466	3473	3486	rVB	2461916	4521443	30.75%	1.701%
99	25.551	3842	3853	3874	rVB	4721748	12842256	87.34%	4.832%
100	26.215	3955	3966	3990	rVB	2179945	6603619	44.91%	2.485%

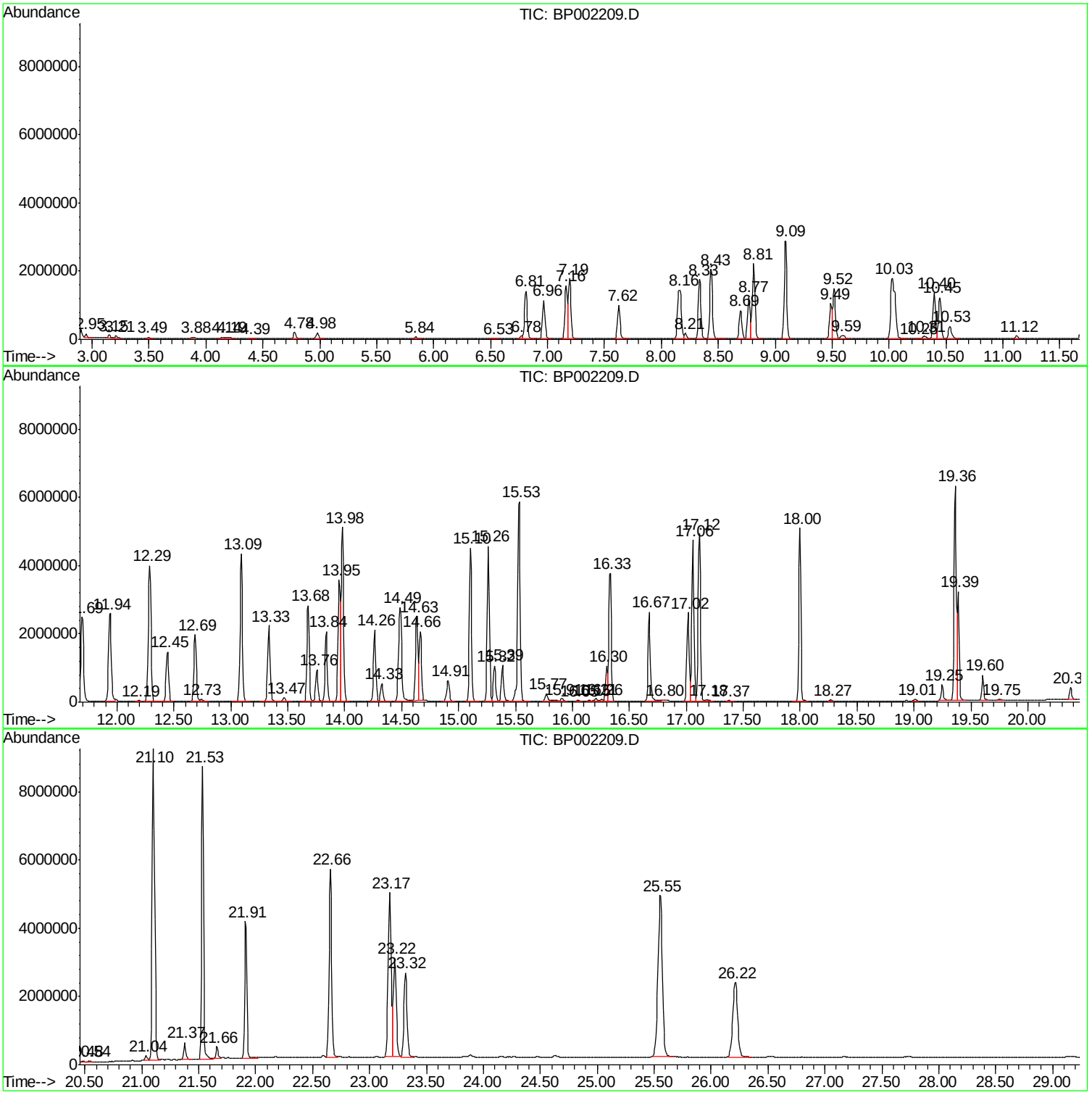
Sum of corrected areas: 265775615

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

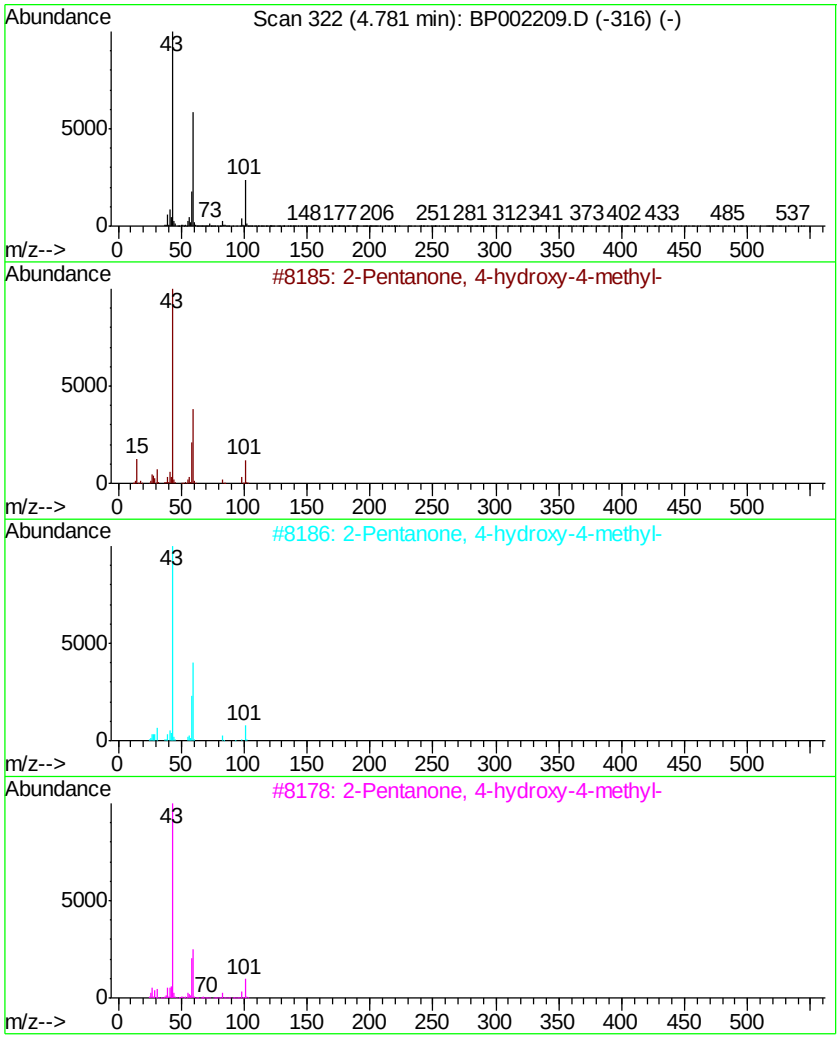
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

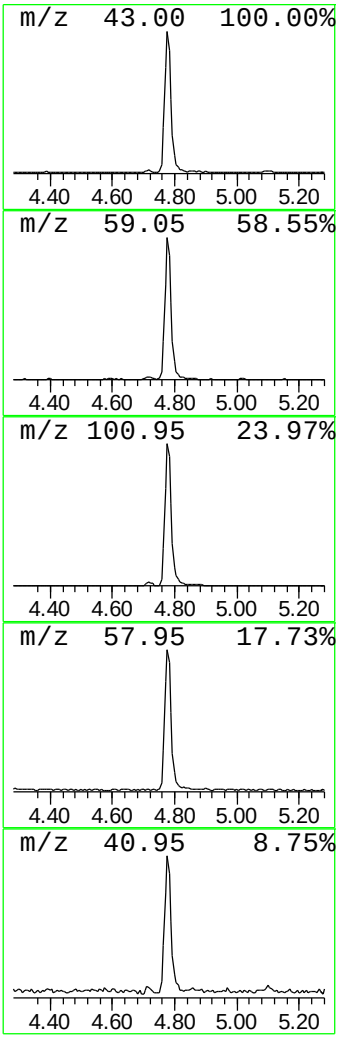
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	3.15 ng/ul	252266	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
5	1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	23





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

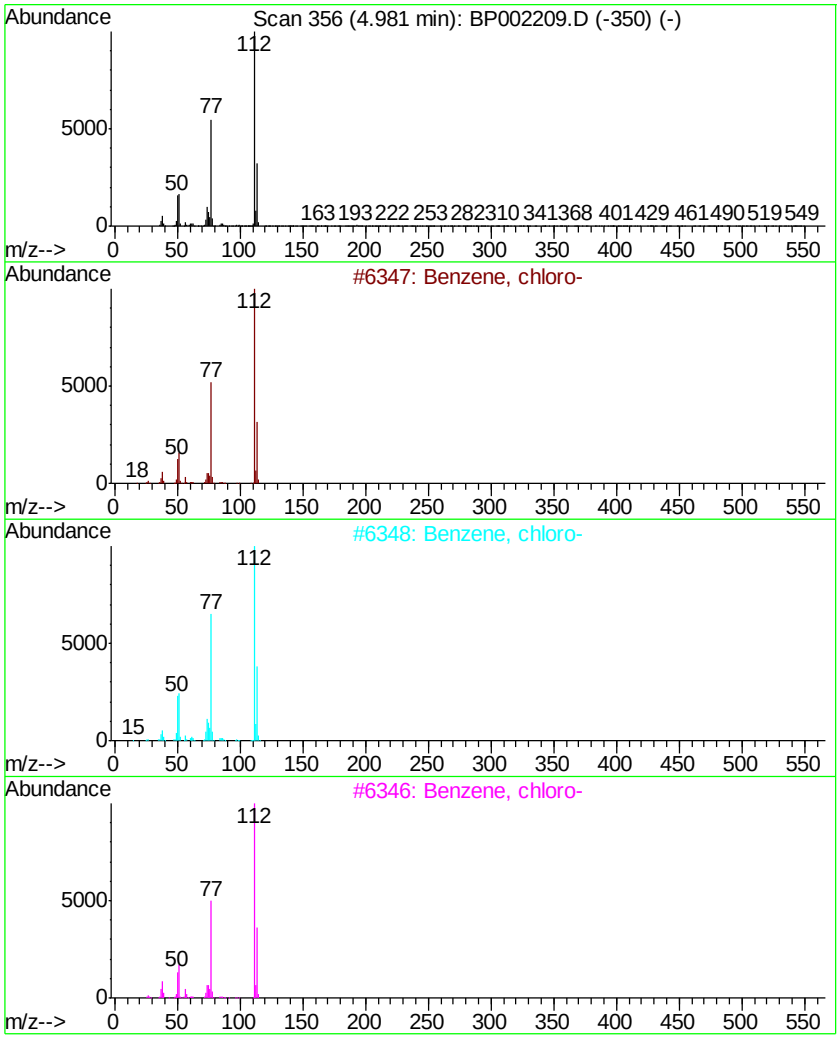
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

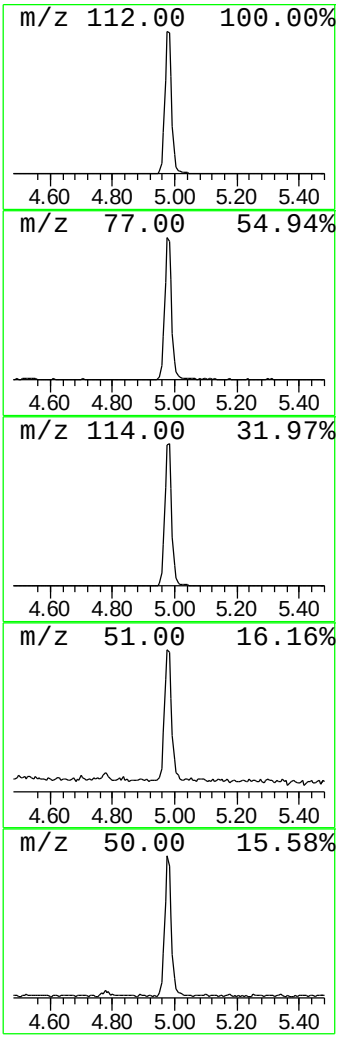
****

Peak Number 2 Benzene, chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	3.29 ng/ul	263629	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Phenol, 4-fluoro-	112	C6H5FO	000371-41-5	23





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

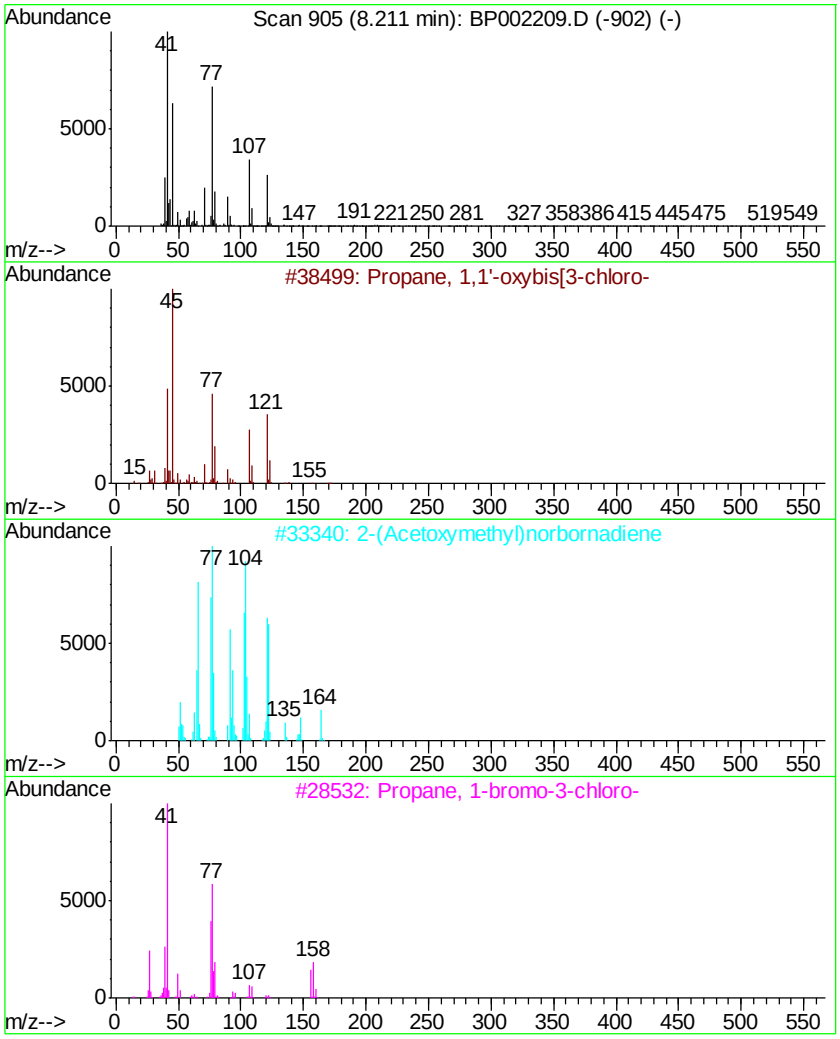
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

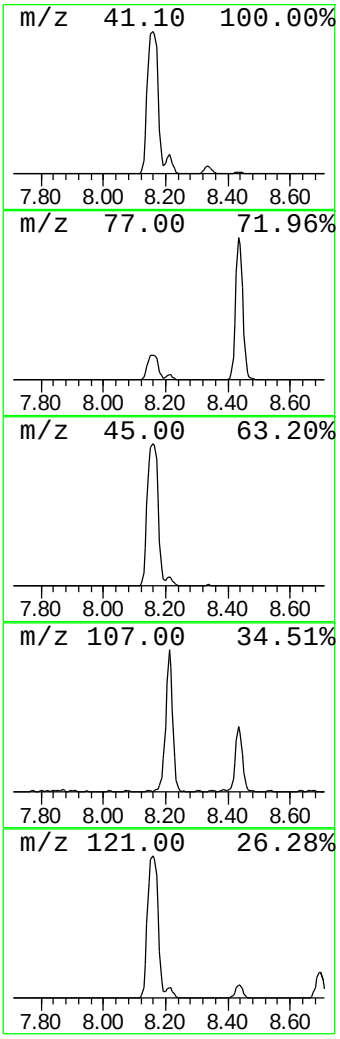
****

Peak Number 3 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.21	2.94 ng/ul	235766	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1'-oxybis[3-chloro-	170	C6H12Cl2O	000629-36-7	40
2	2-(Acetoxymethyl)norbornadiene	164	C10H12O2	056682-74-7	38
3	Propane, 1-bromo-3-chloro-	156	C3H6BrCl	000109-70-6	12
4	Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	10
5	1,3,5-Triazin-2(1H)-one, 4,6-dip...	249	C15H11N3O	001917-44-8	10





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

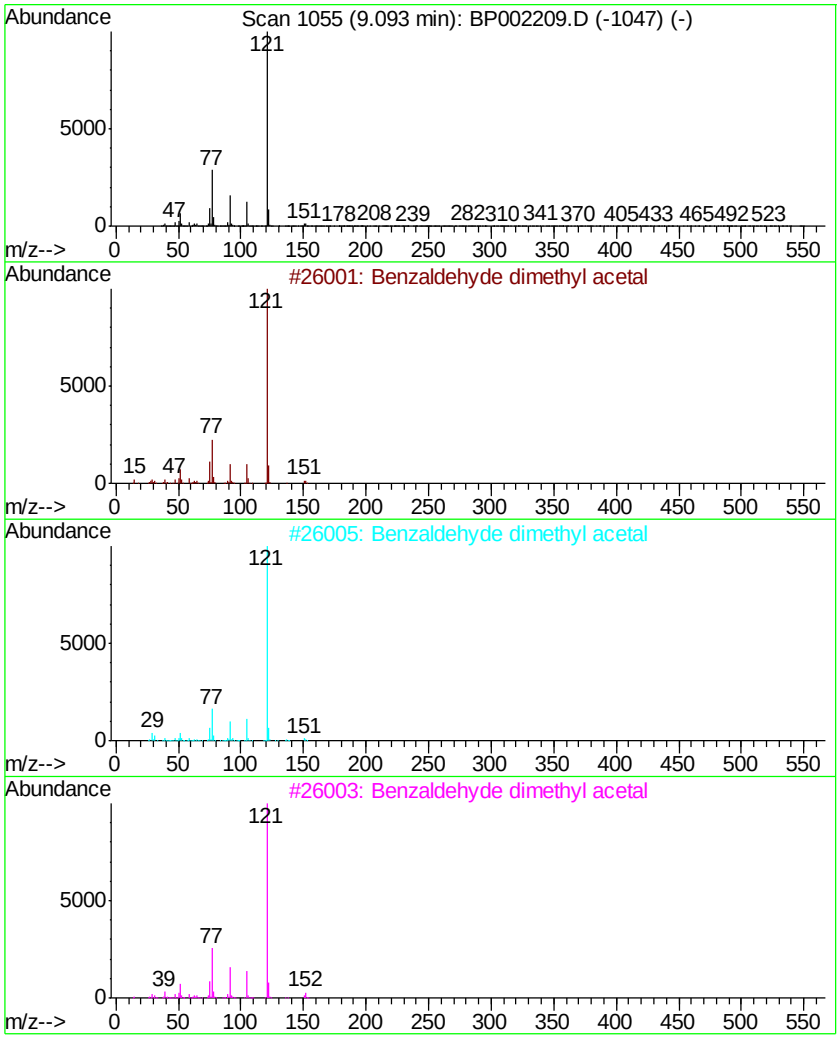
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

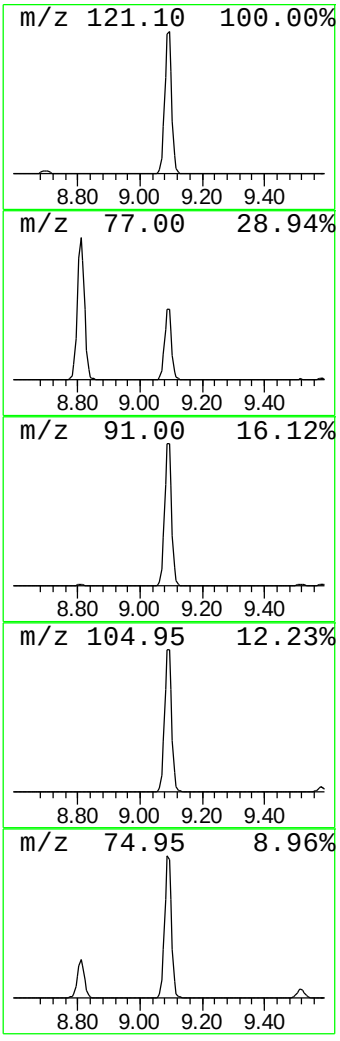
****

Peak Number 4 Benzaldehyde dimethyl acetal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	38.41 ng/ul	4497250	Naphthalene-d8	10.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
2	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
3	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
4	Benzeneacetic acid, .alpha.-meth...	166	C9H10O3	007021-09-2	83
5	Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	78





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

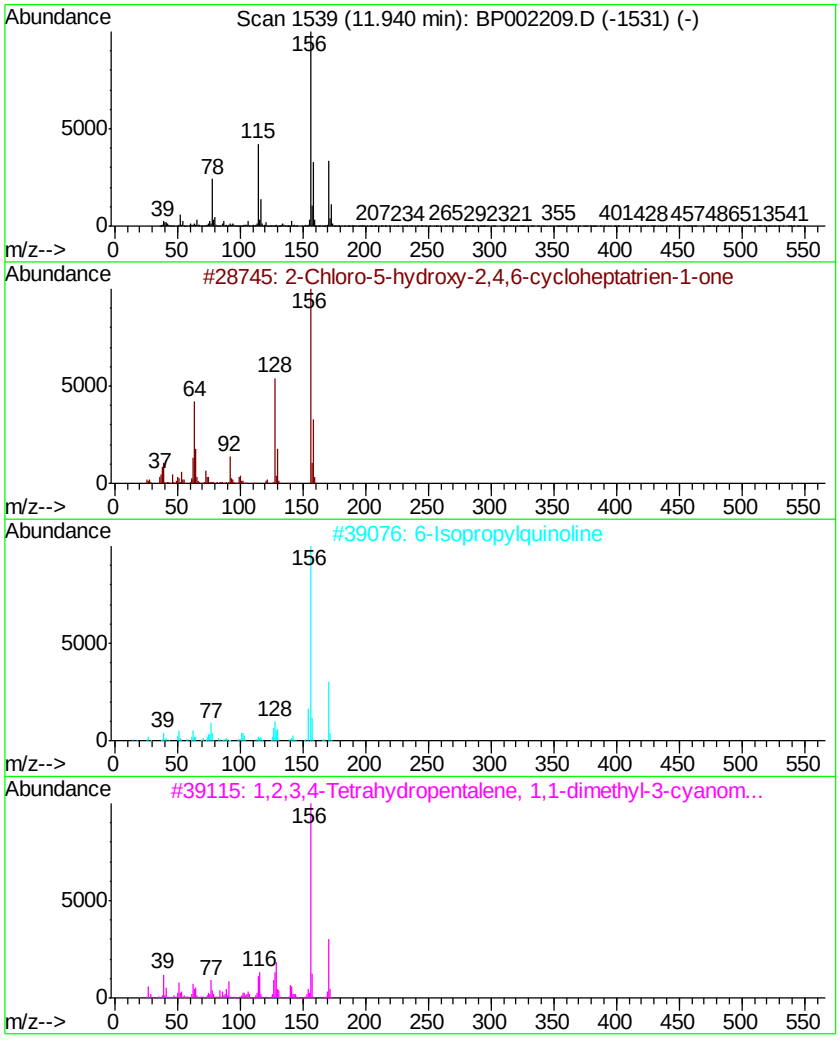
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

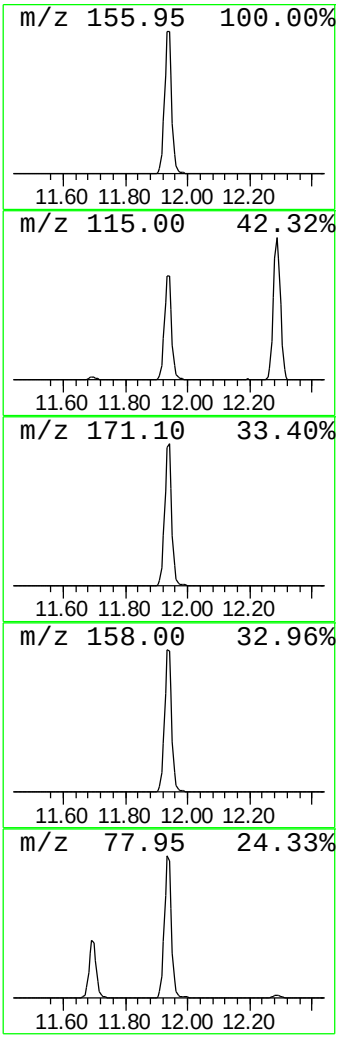
****

Peak Number 5 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	36.74 ng/ul	4301520	Naphthalene-d8	10.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2	6-Isopropylquinoline	171	C12H13N	000135-79-5	38
3	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37
4	1,2,3,3a-Tetrahydropentalene, 1,...	171	C12H13N	1000217-17-0	37
5	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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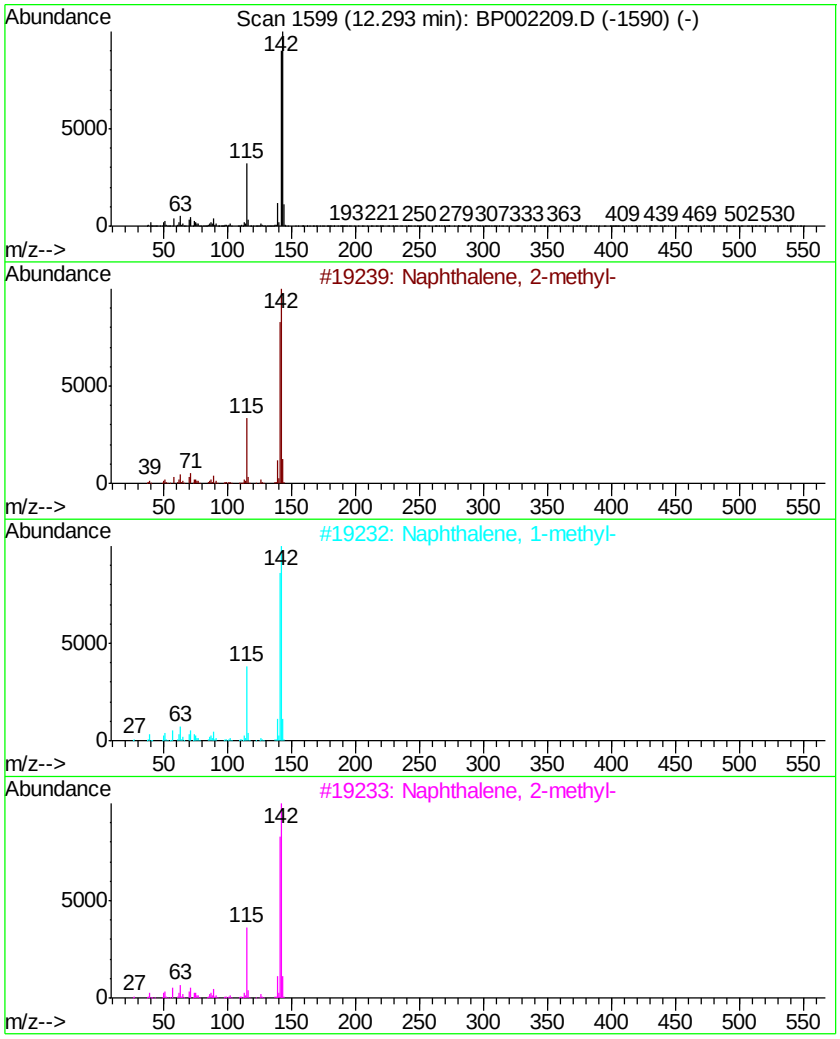
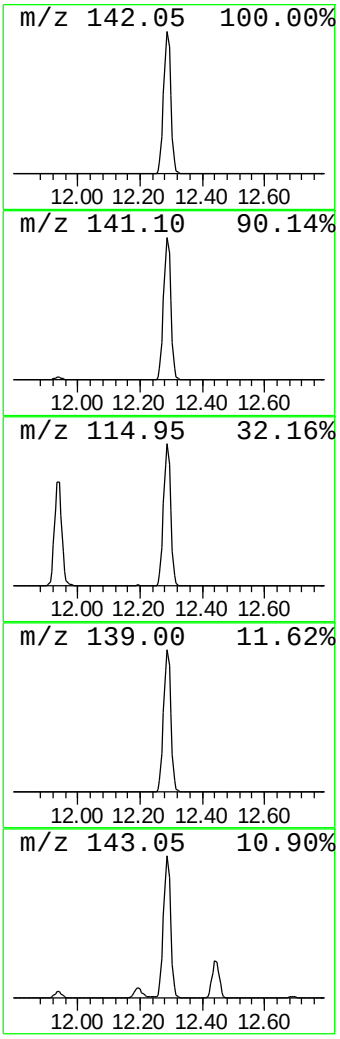
Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	55.12 ng/ul	6454130	Naphthalene-d8	10.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94





Library Search Compound Report

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Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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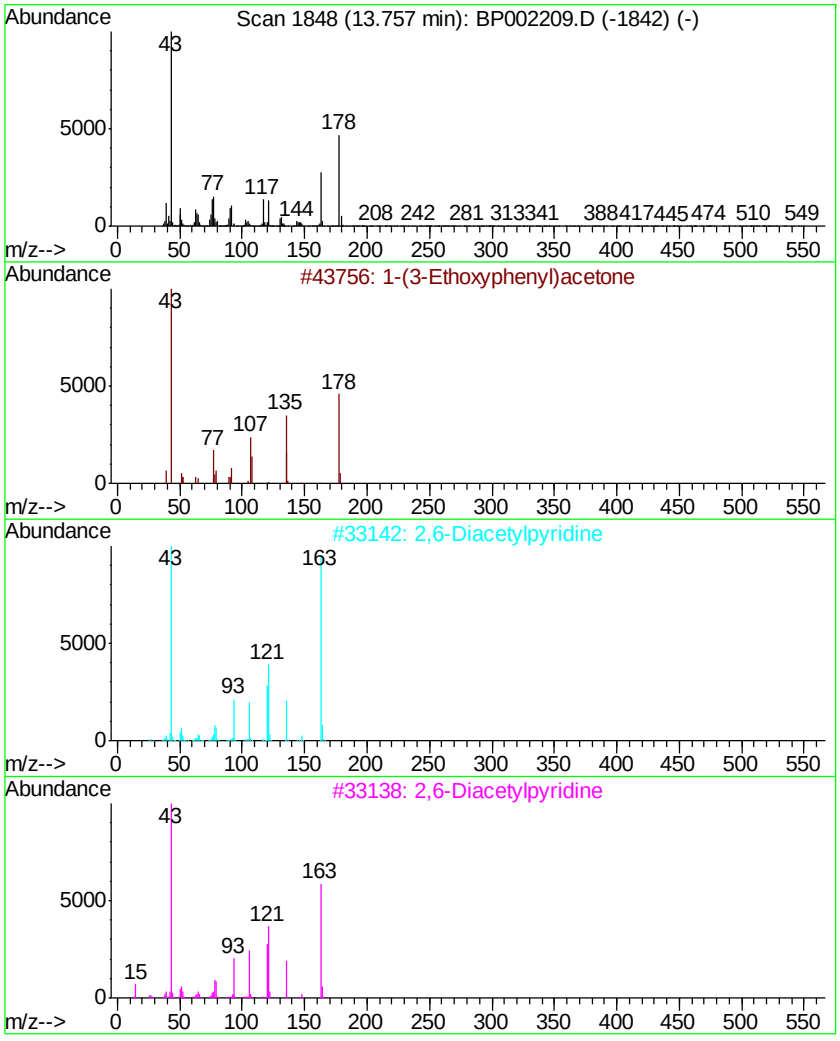
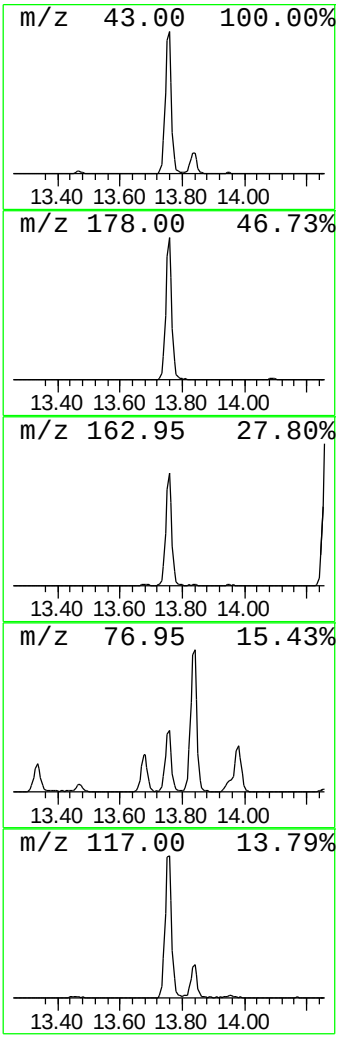
Peak Number 7 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	8.54 ng/ul	1314340	Acenaphthene-d10	14.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	38
2	2,6-Diacetylpyridine	163	C9H9NO2	001129-30-2	10
3	2,6-Diacetylpyridine	163	C9H9NO2	001129-30-2	9
4	Benzeneethanol, .alpha.-methyl-3...	178	C12H18O	054518-11-5	9
5	1-Ethanone, 1-(5-fluoro-1H-1,3-b...	178	C9H7FN2O	1000362-63-4	7





Library Search Compound Report

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Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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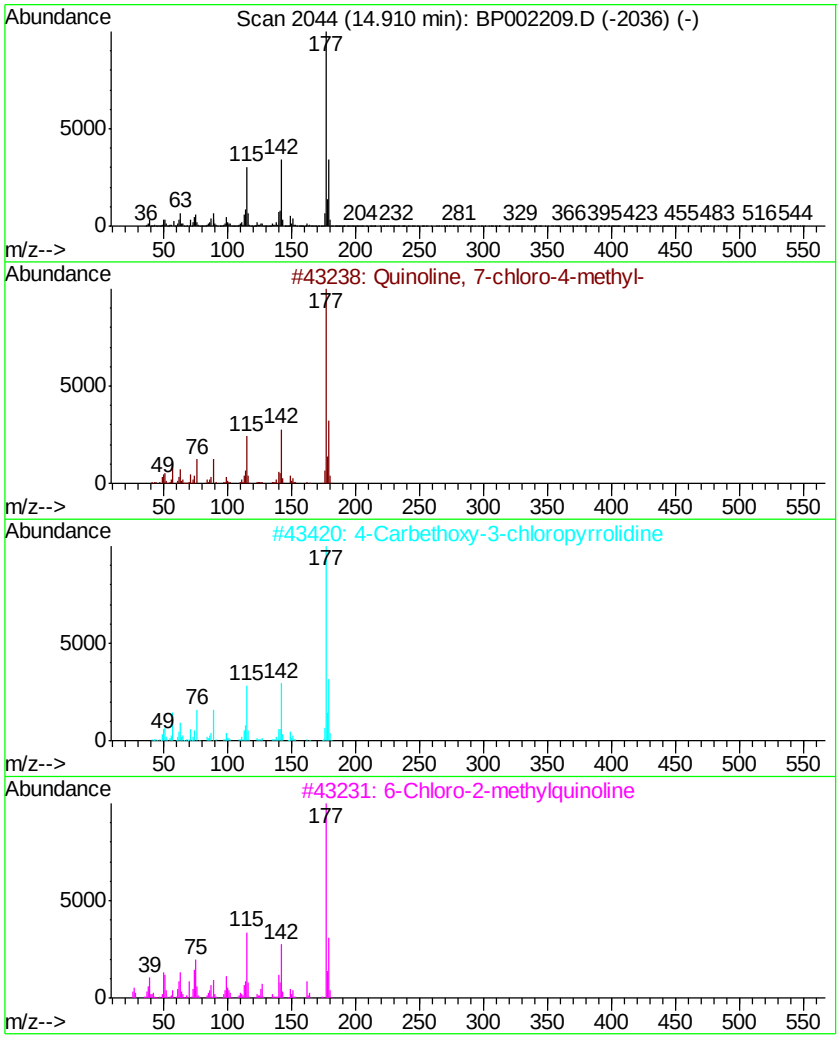
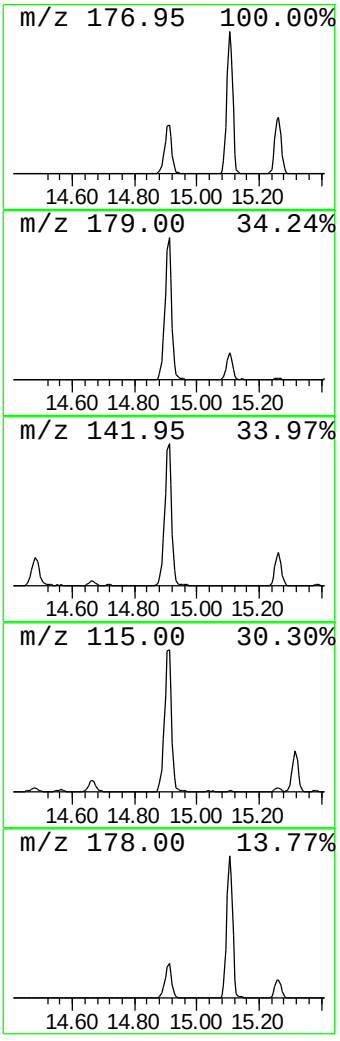
Peak Number 8 Quinoline, 7-chloro-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	6.29 ng/ul	968345	Acenaphthene-d10	14.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline, 7-chloro-4-methyl-	177	C10H8ClN	040941-53-5	94
2	4-Carbethoxy-3-chloropyrrolidine	177	C7H12ClNO2	1000255-97-9	94
3	6-Chloro-2-methylquinoline	177	C10H8ClN	000092-46-6	91
4	Quinoline, 2-chloro-4-methyl-	177	C10H8ClN	000634-47-9	90
5	Quinoline, 7-chloro-2-methyl-	177	C10H8ClN	004965-33-7	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
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Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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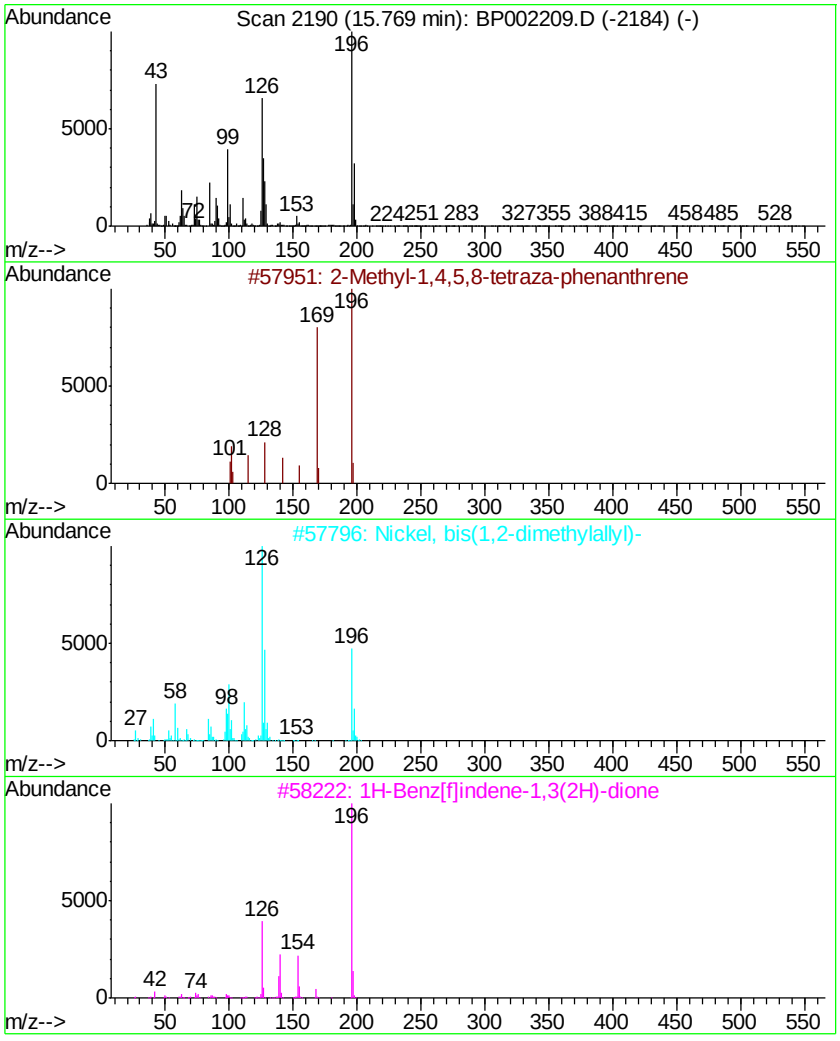
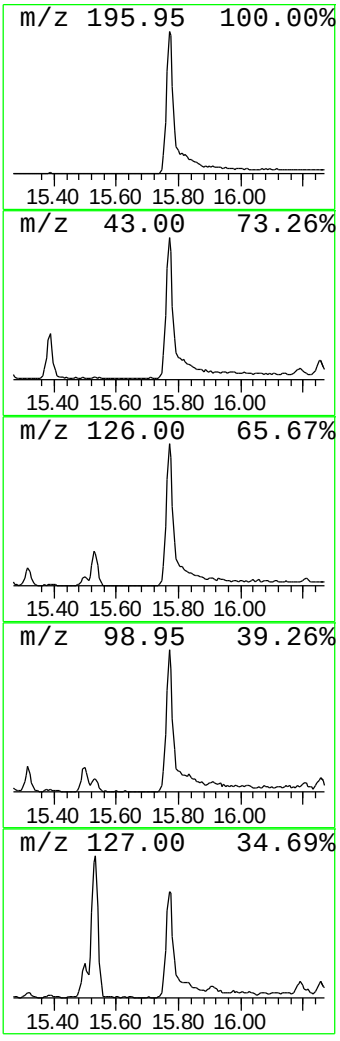
Peak Number 9 unknown-04 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	2.59 ng/ul	480234	Phenanthrene-d10	17.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1,4,5,8-tetraza-phenant...	196	C11H8N4	103538-90-5	38
2	Nickel, bis(1,2-dimethylallyl)-	196	C10H18Ni	1000150-48-3	38
3	1H-Benz[f]indene-1,3(2H)-dione	196	C13H8O2	022734-61-8	32
4	l-Leucine, N-(thiophen-2-carbony...	409	C23H39NO3S	1000325-16-0	30
5	Benzenemethanesulfonamide, 4-chl...	205	C7H8ClNO2S	071799-35-4	27





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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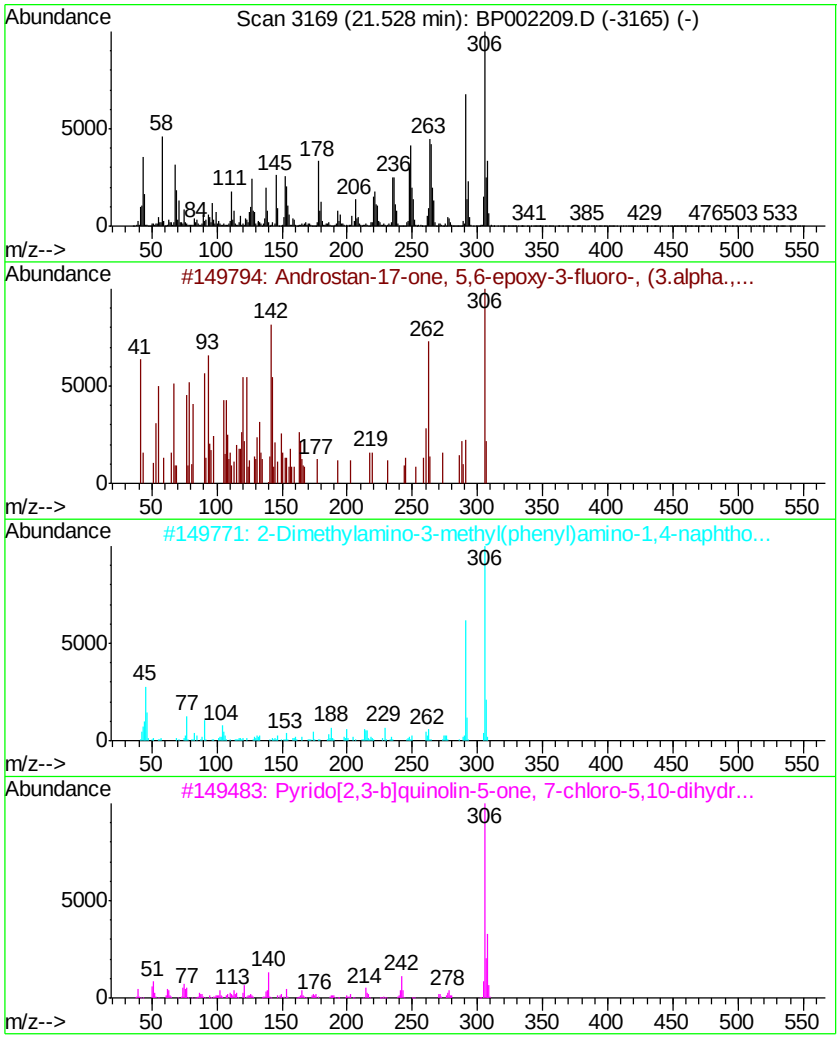
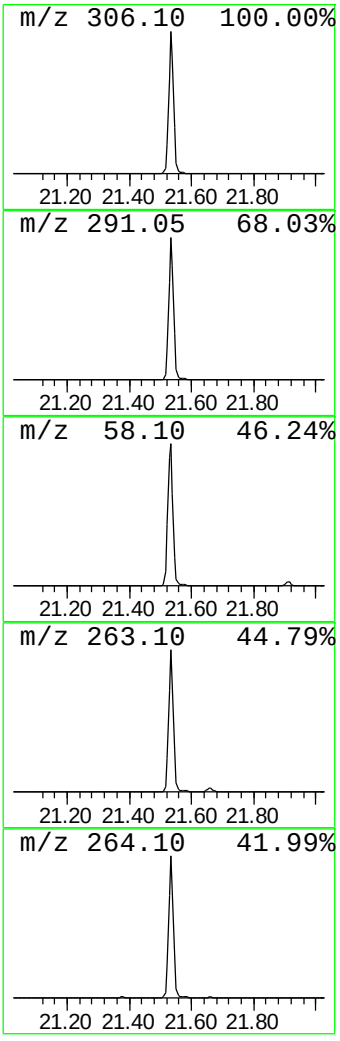
Peak Number 10 Androstan-17-one, 5,6-epoxy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.53	13.42 ng/ul	9863510	Chrysene-d12	21.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Androstan-17-one, 5,6-epoxy-3-fl...	306	C19H27FO2	028344-36-7	92
2	2-Dimethylamino-3-methyl(phenyl)...	306	C19H18N2O2	134614-10-1	42
3	Pyrido[2,3-b]quinolin-5-one, 7-c...	306	C18H11ClN2O	069750-07-8	38
4	Acetamide, 2-[(4,5-dihydro-6-met...	306	C13H14N4O3S	1000317-32-8	25
5	1,1':2',1''-Terphenyl, 4'-phenyl-	306	C24H18	001165-53-3	22





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031220\
Data File : BP002209.D
Acq On : 13 Mar 2020 10:31
Operator : CG/JU
Sample : L1840-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBER9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.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
2-Pentanone, 4-hy...	4.78	3.1	ng/ul	252266	1	7.62	1602470	20.0
Benzene, chloro-	4.98	3.3	ng/ul	263629	1	7.62	1602470	20.0
unknown-01	8.21	2.9	ng/ul	235766	1	7.62	1602470	20.0
Benzaldehyde dime...	9.09	38.4	ng/ul	4497250	2	10.39	2341840	20.0
unknown-02	11.94	36.7	ng/ul	4301520	2	10.39	2341840	20.0
Naphthalene, 1-me...	12.29	55.1	ng/ul	6454130	2	10.39	2341840	20.0
unknown-03	13.76	8.5	ng/ul	1314340	3	14.26	3079660	20.0
Quinoline, 7-chlo...	14.91	6.3	ng/ul	968345	3	14.26	3079660	20.0
unknown-04	15.77	2.6	ng/ul	480234	4	17.02	3713860	20.0
Androstan-17-one,...	21.53	13.4	ng/ul	9863510	5	21.12	14704300	20.0