LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM026531.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.263	59	64	74	rVB	23637	41608	1.85%	0.236%
2	3.663	128	132	138	rVB	19103	27961	1.24%	0.158%
3	4.093	200	205	212	rVB8	7118	13861	0.61%	0.078%
4	4.346	243	248	254	rBV2	16079	26820	1.19%	0.152%
5	4.481	267	271	276	rBV5	5875	11082	0.49%	0.063%
6	4.957	346	352	355	rBV6	3562	7909	0.35%	0.045%
7	5.287	402	408	416	rBV2	14312	26206	1.16%	0.148%
8	5.969	519	524	530	rVV8	7278	13537	0.60%	0.077%
9	6.081	538	543	548	rVB5	9901	14565	0.65%	0.082%
10	6.375	589	593	600	rVB5	10949	17206	0.76%	0.097%
11	6.469	600	609	613	rBV5	11814	22161	0.98%	0.126%
12	6.616	625	634	640	rBV3	18462	46149	2.05%	0.261%
13	6.740	650	655	661	rBV	25860	44872	1.99%	0.254%
14	6.793	661	664	672	rVV3	14104	27110	1.20%	0.154%
15	6.922	680	686	691	rVV2	32951	54316	2.41%	0.308%
16	6.963	691	693	699	rVB5	5302	9298	0.41%	0.053%
17	7.216	732	736	740	rVB4	6006	7843	0.35%	0.044%
18	7.375	755	763	771	rBV	427070	665753	29.53%	3.770%
19	7.516	780	787	794	rVV	210220	302908	13.43%	1.715%
20	7.681	812	815	822	rVB7	6921	9396	0.42%	0.053%
21	7.934	852	858	867	rBV2	20389	38324	1.70%	0.217%
22	8.046	870	877	881	rBV5	3847	8627	0.38%	0.049%
23	8.104	881	887	890	rBV2	30853	46919	2.08%	0.266%
24	8.157	890	896	911	rVB	1423316	2254740	100.00%	12.769%
25	8.363	927	931	934	rVB3	7387	8965	0.40%	0.051%
26	8.528	953	959	967	rVV	38893	69815	3.10%	0.395%
27	8.604	967	972	978	rVB	135652	201195	8.92%	1.139%
28	8.757	994	998	1002	rVB4	6766	8308	0.37%	0.047%
29	8.940	1024	1029	1035	rBV7	3542	8857	0.39%	0.050%
30	9.240	1074	1080	1085	rBV	30291	54798	2.43%	0.310%
31	9.522	1119	1128	1130	rVV	89654	150534	6.68%	0.853%
32	9.557	1130	1134	1141	rVV	258647	407071	18.05%	2.305%
33	9.687	1150	1156	1162	rVV2	24109	45542	2.02%	0.258%
34	9.792	1168	1174	1186	rVB3	41611	94502	4.19%	0.535%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Title : SVOA CALIBRATION

35	10.016	1205	1212	1221	rVB	86097	142928	6.34%	0.809%
36	10.134	1226	1232	1239	rVV	643928	1038703	46.07%	5.882%
37	10.210	1239	1245	1250	rVV2	144457	244850	10.86%	1.387%
38	10.257	1250	1253	1255	rVV4	13334	22285	0.99%	0.126%
39	10.298	1255	1260	1273	rVV2	37169	92752	4.11%	0.525%
40	10.398	1274	1277	1282	rVV6	5081	9161	0.41%	0.052%
41	10.498	1289	1294	1298	rVV2	32119	50739	2.25%	0.287%
42	10.581	1305	1308	1314	rVB6	4926	9457	0.42%	0.054%
43	10.904	1356	1363	1366	rBV7	4974	9359	0.42%	0.053%
44	10.992	1373	1378	1386	rBV3	46856	93170	4.13%	0.528%
45	11.063	1386	1390	1398	rVB	12141	21101	0.94%	0.119%
46	11.139	1398	1403	1409	rVB6	4615	8277	0.37%	0.047%
47	11.510	1458	1466	1472	rVB5	6733	14225	0.63%	0.081%
48	11.669	1487	1493	1497	rVB7	5630	9333	0.41%	0.053%
49	11.904	1525	1533	1538	rVB3	21757	38229	1.70%	0.216%
50	12.016	1544	1552	1564	rBV	110514	223227	9.90%	1.264%
51	12.181	1575	1580	1585	rBV7	7936	12603	0.56%	0.071%
52	12.339	1602	1607	1612	rBV2	26200	41701	1.85%	0.236%
53	12.486	1625	1632	1641	rBV	159930	235882	10.46%	1.336%
54	12.680	1661	1665	1670	rBV5	7210	10230	0.45%	0.058%
55	12.751	1673	1677	1681	rVB4	7096	10762	0.48%	0.061%
56	12.886	1694	1700	1706	rBV	52568	82250	3.65%	0.466%
57	12.951	1706	1711	1722	rVB2	35619	68826	3.05%	0.390%
58	13.086	1728	1734	1740	rVB9	6371	11087	0.49%	0.063%
59	13.463	1793	1798	1802	rBV	98110	137876	6.11%	0.781%
60	13.510	1802	1806	1815	rVB	41520	69584	3.09%	0.394%
61	13.722	1836	1842	1848	rVV	108375	167616	7.43%	0.949%
62	13.869	1862	1867	1879	rBV	75726	149169	6.62%	0.845%
63	13.963	1879	1883	1887	rVV5	9897	16239	0.72%	0.092%
64	14.033	1889	1895	1903	rVV	1199488	1724135	76.47%	9.764%
65	14.092	1903	1905	1912	rVB2	15210	23230	1.03%	0.132%
66	14.180	1918	1920	1928	rVB8	5200	11311	0.50%	0.064%
67	14.351	1943	1949	1954	rBV9	5731	13882	0.62%	0.079%
68	14.504	1969	1975	1981	rBV8	11131	20444	0.91%	0.116%
69	14.580	1983	1988	1996	rVB10	8334	17679	0.78%	0.100%
70	14.663	1996	2002	2005	rBV6	5397	8831	0.39%	0.050%
71	14.810	2018	2027	2035	rBV2	68106	161367	7.16%	0.914%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Title : SVOA CALIBRATION

72	14.974	2051	2055	2060	rBV3	9846	15348	0.68%	0.087%
73	15.039	2060	2066	2071	rBV	141987	194043	8.61%	1.099%
74	15.098	2072	2076	2081	rVB2	30923	42875	1.90%	0.243%
75	15.192	2086	2092	2099	rBV4	29707	49107	2.18%	0.278%
76	15.304	2108	2111	2116	rVB6	5997	9399	0.42%	0.053%
77	15.421	2126	2131	2133	rBV4	9870	14449	0.64%	0.082%
78	15.451	2133	2136	2143	rVB5	14640	20543	0.91%	0.116%
79	15.569	2152	2156	2164	rVV	61521	91121	4.04%	0.516%
80	15.927	2211	2217	2225	rBV	11693	31895	1.41%	0.181%
81	16.086	2240	2244	2251	rVV	26779	41635	1.85%	0.236%
82	16.151	2251	2255	2258	rVB4	6586	8504	0.38%	0.048%
83	16.245	2266	2271	2274	rBV6	11351	16499	0.73%	0.093%
84	16.574	2324	2327	2333	rBV6	5209	10350	0.46%	0.059%
85	16.792	2357	2364	2370	rBV	1325387	1832815	81.29%	10.380%
86	16.892	2375	2381	2390	rVB	130089	190362	8.44%	1.078%
87	16.998	2394	2399	2405	rBV	264538	347652	15.42%	1.969%
88	17.210	2432	2435	2439	rBV5	10722	15558	0.69%	0.088%
89	17.727	2520	2523	2528	rBV4	13691	17377	0.77%	0.098%
90	18.527	2654	2659	2663	rBV8	9236	14703	0.65%	0.083%
91	18.592	2668	2670	2676	rBV7	9009	15377	0.68%	0.087%
92	18.862	2713	2716	2721	rBV2	10777	13691	0.61%	0.078%
93	19.098	2752	2756	2767	rBV2	27701	57798	2.56%	0.327%
94	19.198	2768	2773	2781	rVV	112782	164917	7.31%	0.934%
95	20.756	3035	3038	3042	rBV2	51111	59739	2.65%	0.338%
96	21.003	3075	3080	3091	rBV	1827724	2211836	98.10%	12.526%
97	22.509	3333	3336	3341	rVB2	96685	118528	5.26%	0.671%
98	23.027	3421	3424	3428	rVB2	136648	173344	7.69%	0.982%
99	23.092	3429	3435	3445	rVB	1129150	1871690	83.01%	10.600%
100	23.609	3520	3523	3529	rVB2	124269	195418	8.67%	1.107%

Sum of corrected areas: 17657831

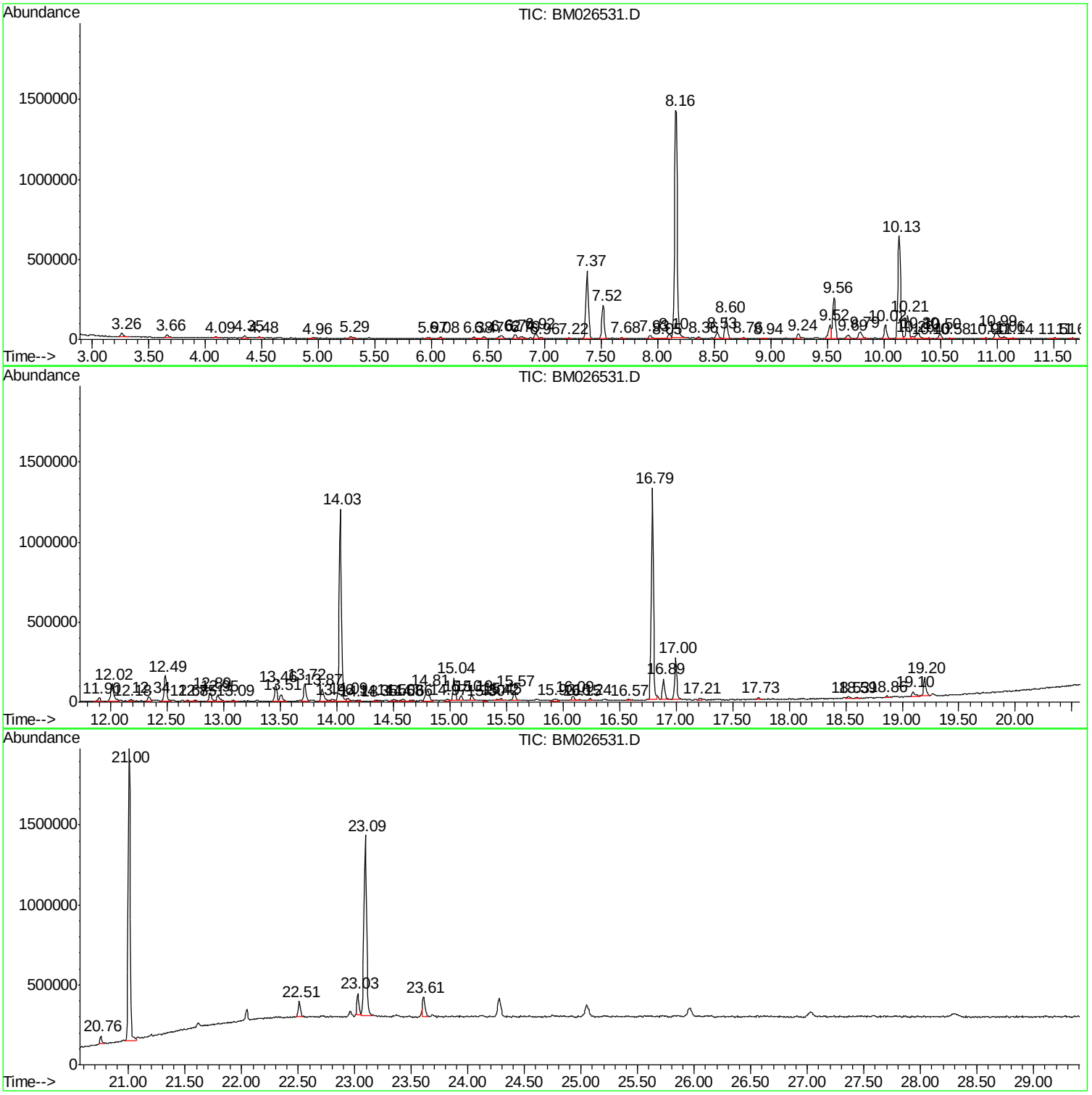
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

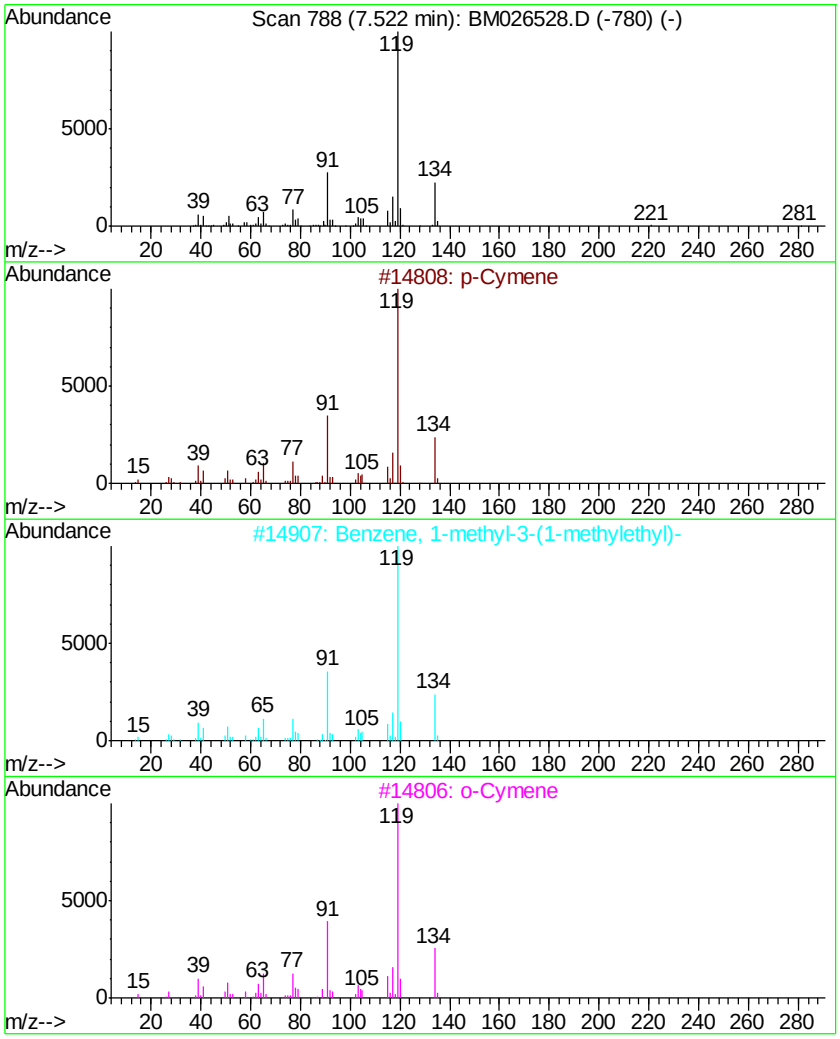
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

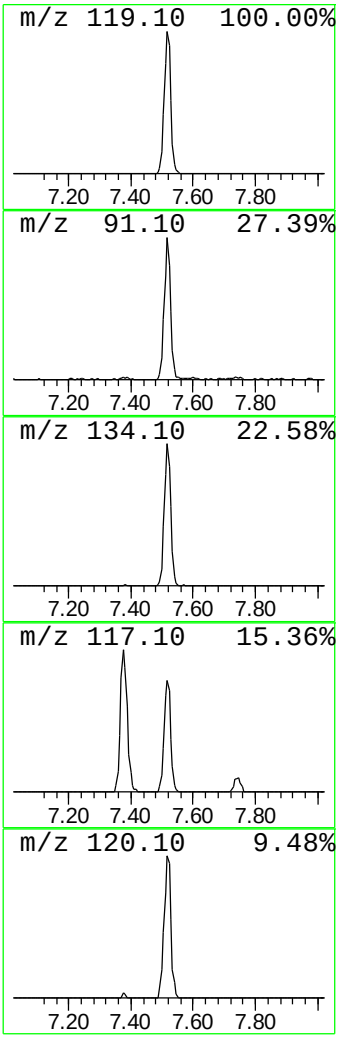
****

Peak Number 1 p-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	9.10 ng/ul	302908	1,4-Dichlorobenzene-d4	7.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	96
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

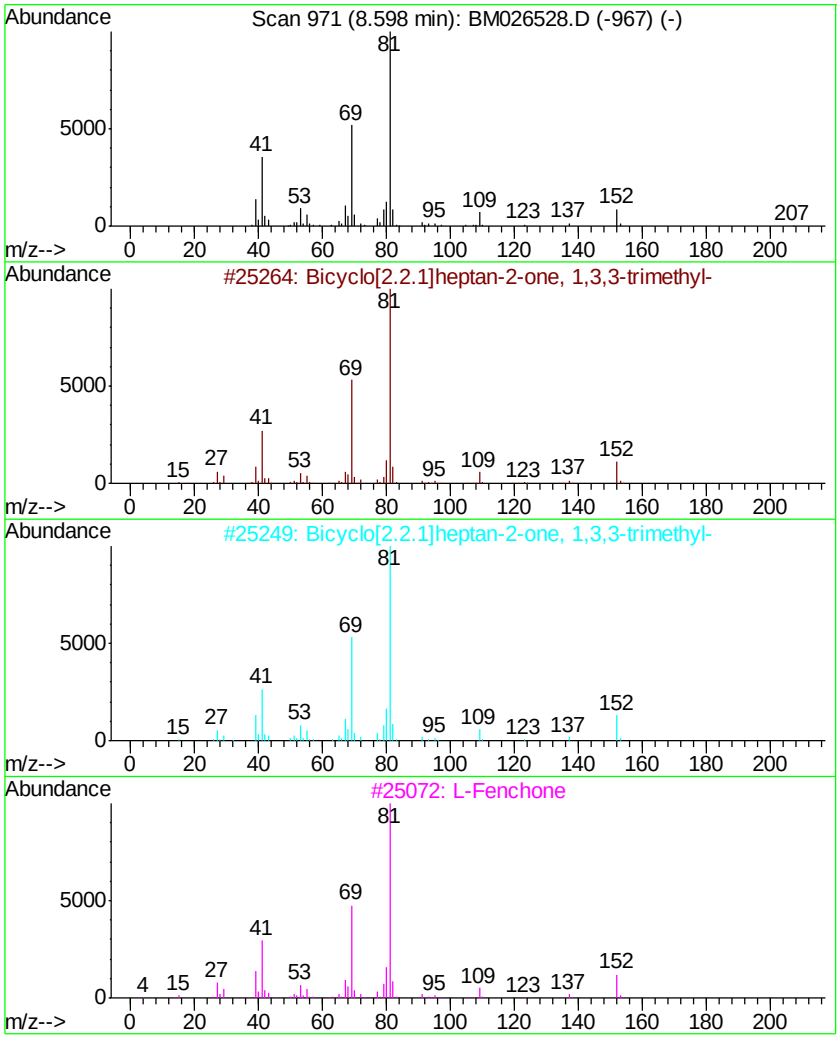
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

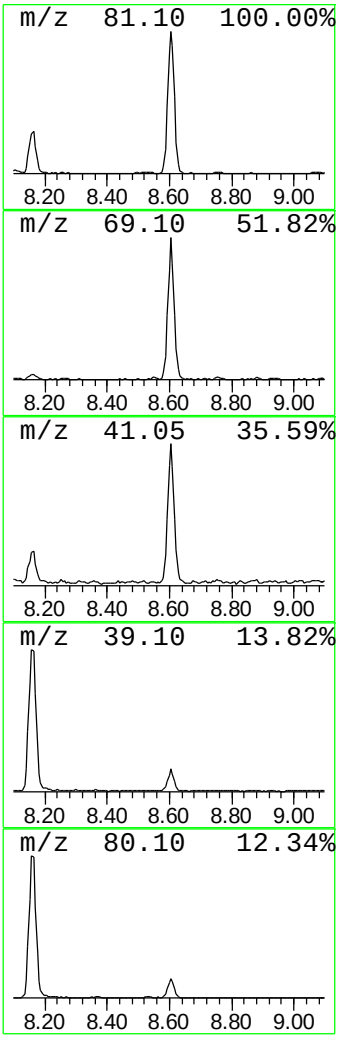
****

Peak Number 2 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	6.04 ng/ul	201195	1,4-Dichlorobenzene-d4	7.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
3	L-Fenchone	152	C10H16O	007787-20-4	93
4	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	90
5	D-Fenchone	152	C10H16O	004695-62-9	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

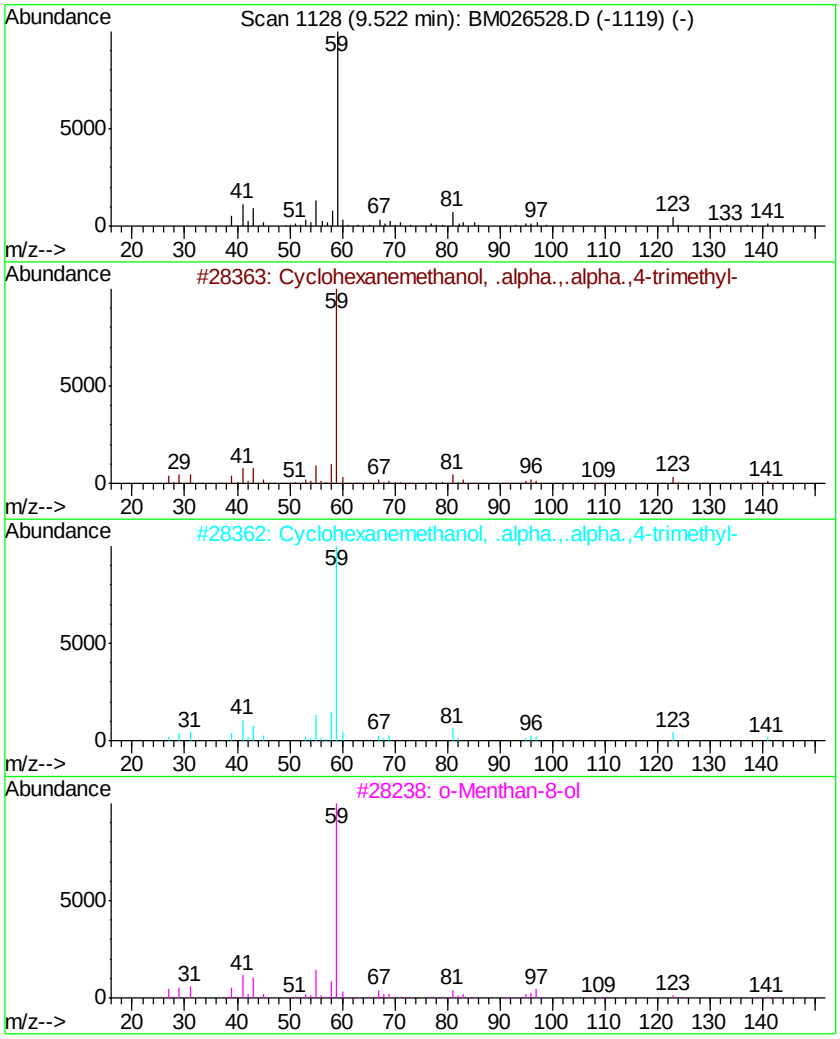
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

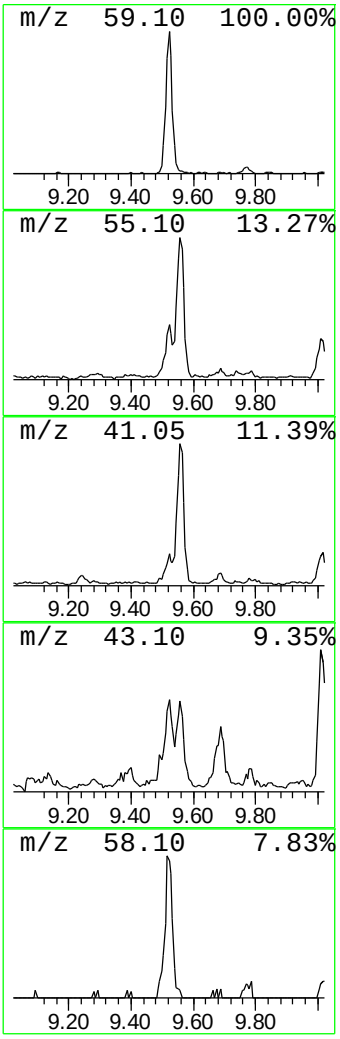
****

Peak Number 3 Cyclohexanemethanol, .alpha... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.52	2.90 ng/ul	150534	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	78
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	74
3	o-Menthan-8-ol	156	C10H20O	343855-44-7	72
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	56
5	2,3-Dimethyl-4-penten-2-ol	114	C7H14O	019781-52-3	56





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

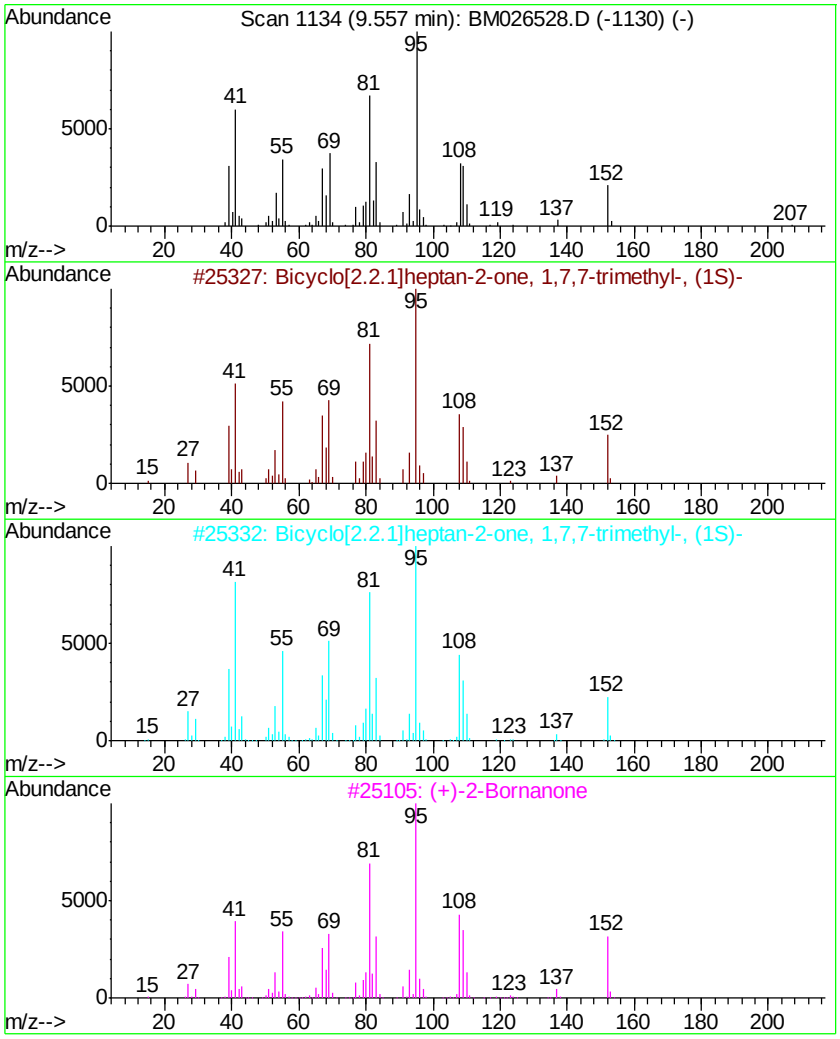
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

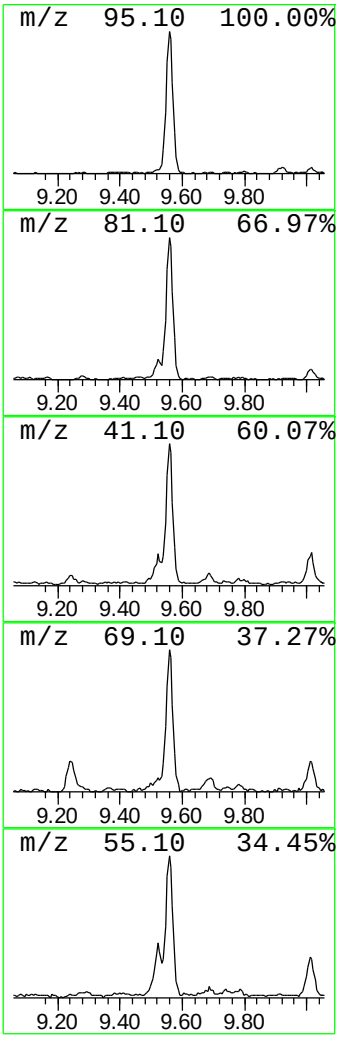
****

Peak Number 4 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	7.84 ng/ul	407071	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
2	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
4	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	96





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

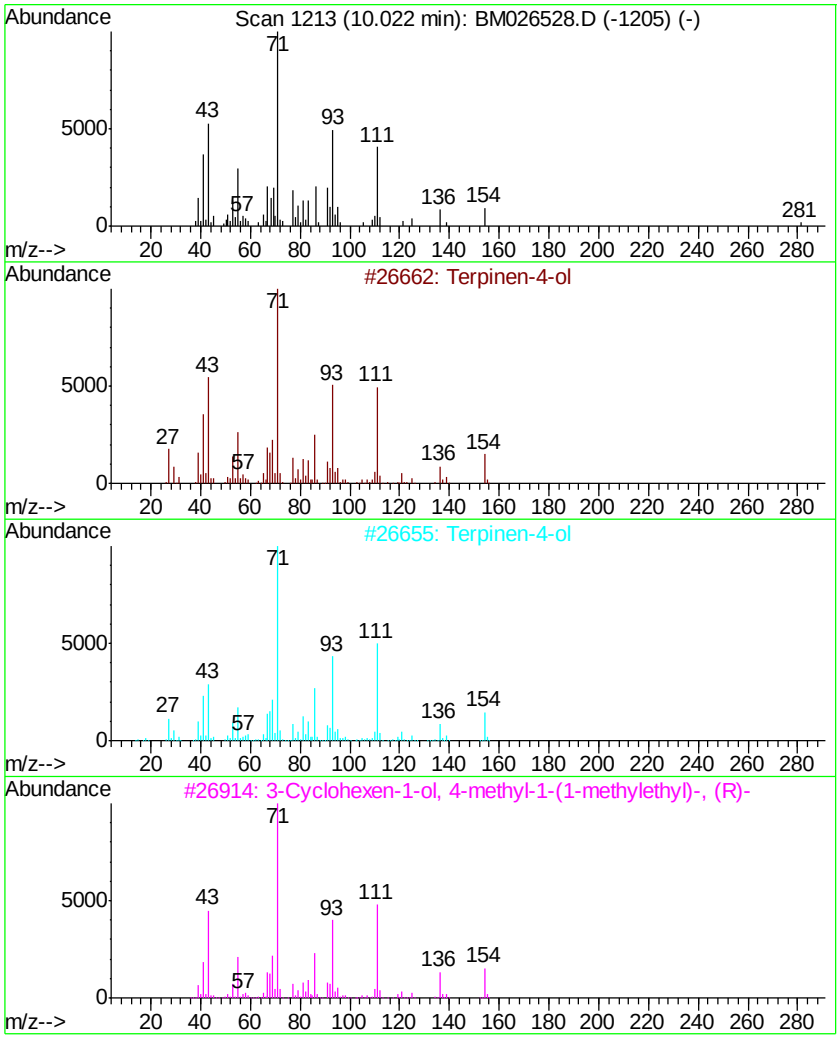
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

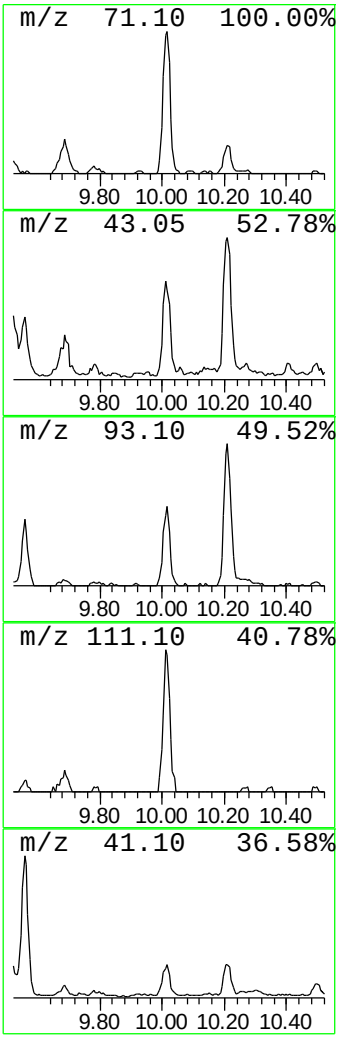
****

Peak Number 5 Terpinen-4-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	2.75 ng/ul	142928	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Terpinen-4-ol	154	C10H18O	000562-74-3	91
2	Terpinen-4-ol	154	C10H18O	000562-74-3	91
3	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	81
4	Terpinen-4-ol	154	C10H18O	000562-74-3	64
5	Terpinen-4-ol	154	C10H18O	000562-74-3	58





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

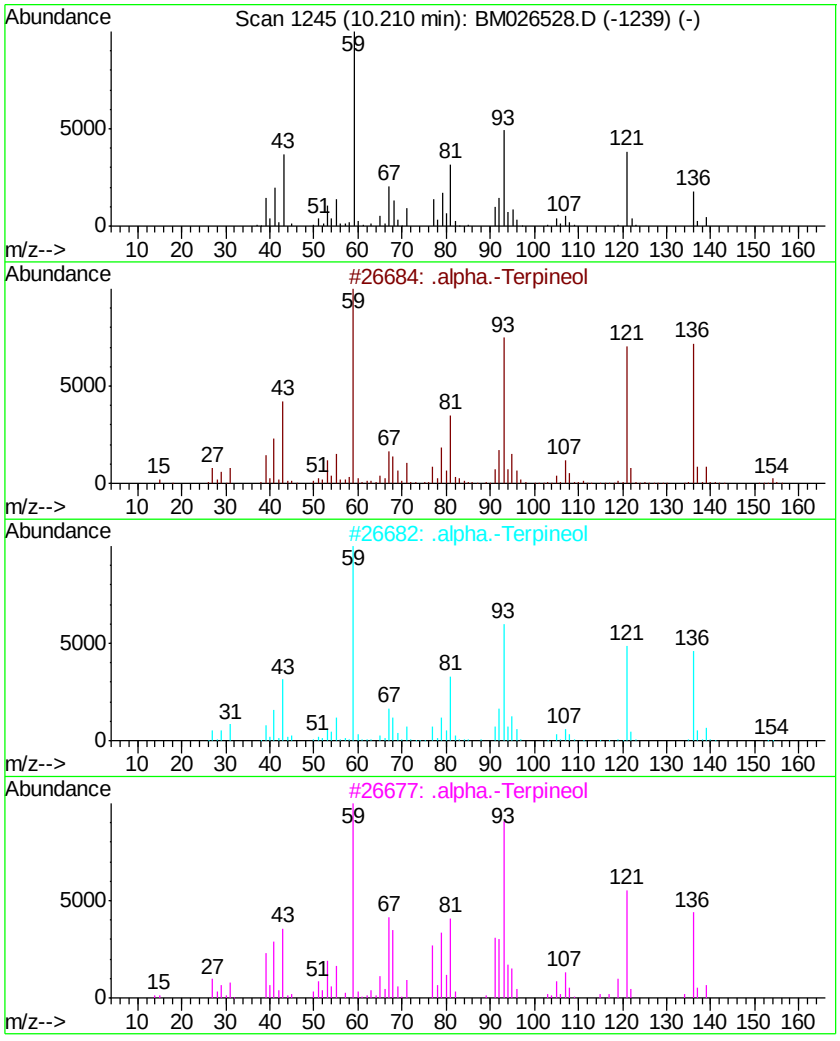
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

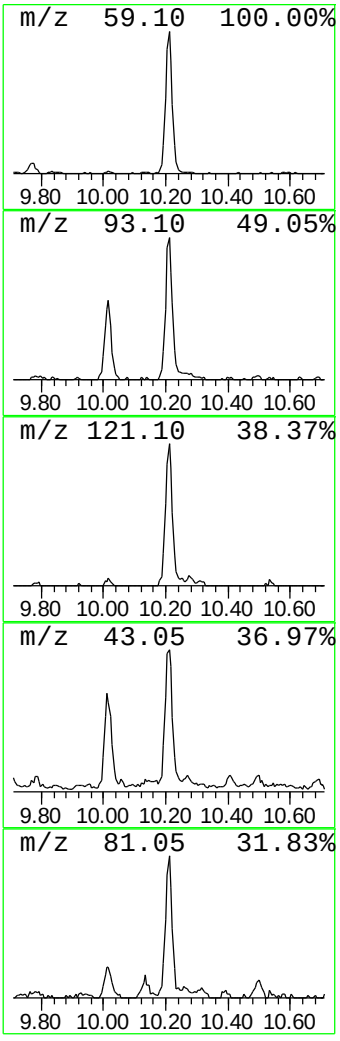
****

Peak Number 6 .alpha.-Terpineol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.21	4.71 ng/ul	244850	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Terpineol	154	C10H18O	000098-55-5	72
2	.alpha.-Terpineol	154	C10H18O	000098-55-5	72
3	.alpha.-Terpineol	154	C10H18O	000098-55-5	46
4	.alpha.-Terpineol	154	C10H18O	000098-55-5	43
5	L-.alpha.-Terpineol	154	C10H18O	010482-56-1	43





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Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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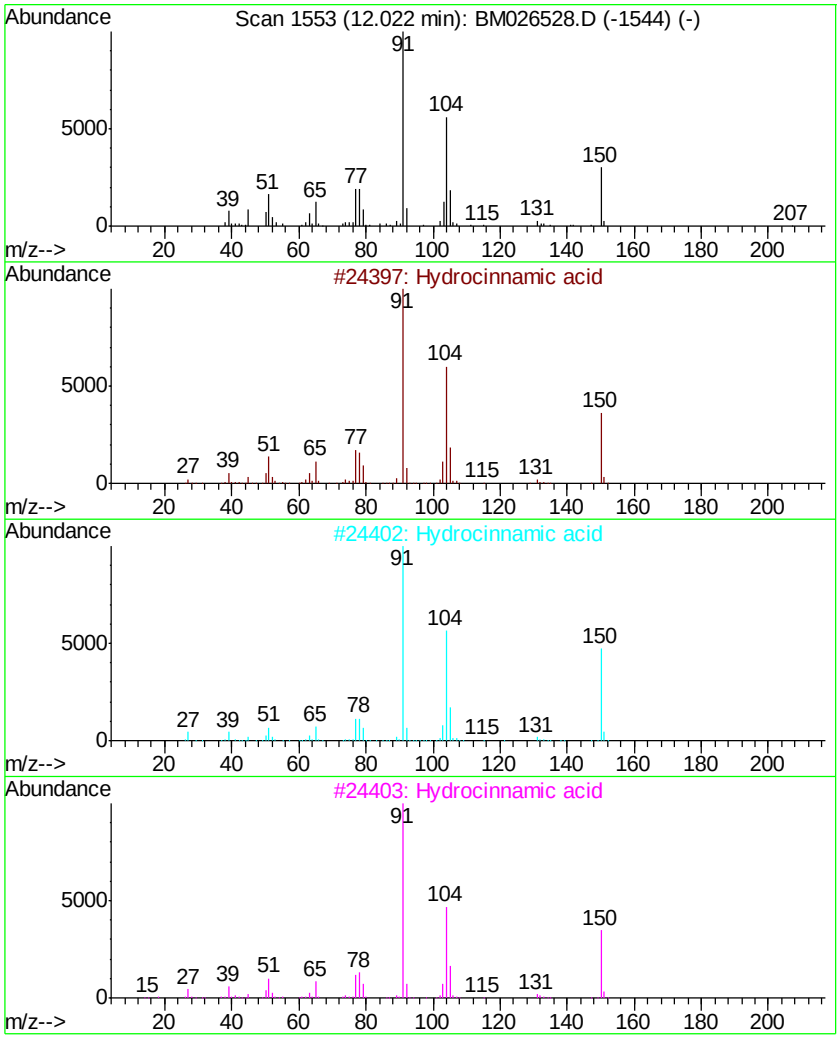
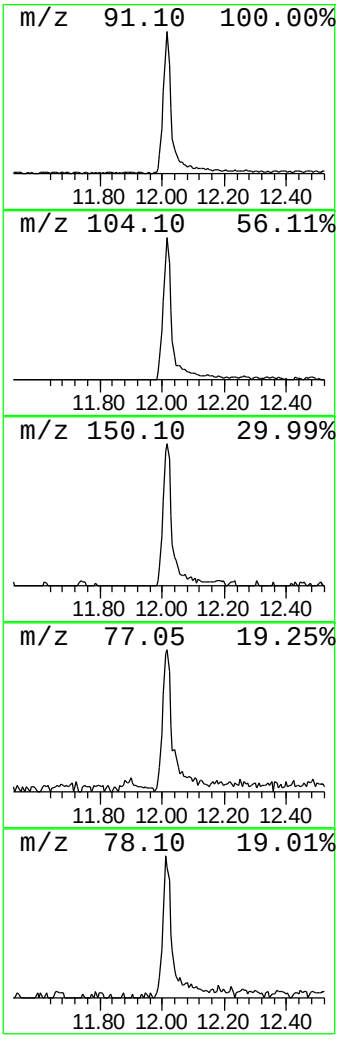
Peak Number 7 Hydrocinnamic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	4.30 ng/ul	223227	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrocinnamic acid	150	C9H10O2	000501-52-0	97
2	Hydrocinnamic acid	150	C9H10O2	000501-52-0	91
3	Hydrocinnamic acid	150	C9H10O2	000501-52-0	90
4	4-Nitrosophenyl-.beta.-phenylpro...	255	C15H13NO3	1000129-40-0	90
5	Hydrocinnamic acid	150	C9H10O2	000501-52-0	68





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Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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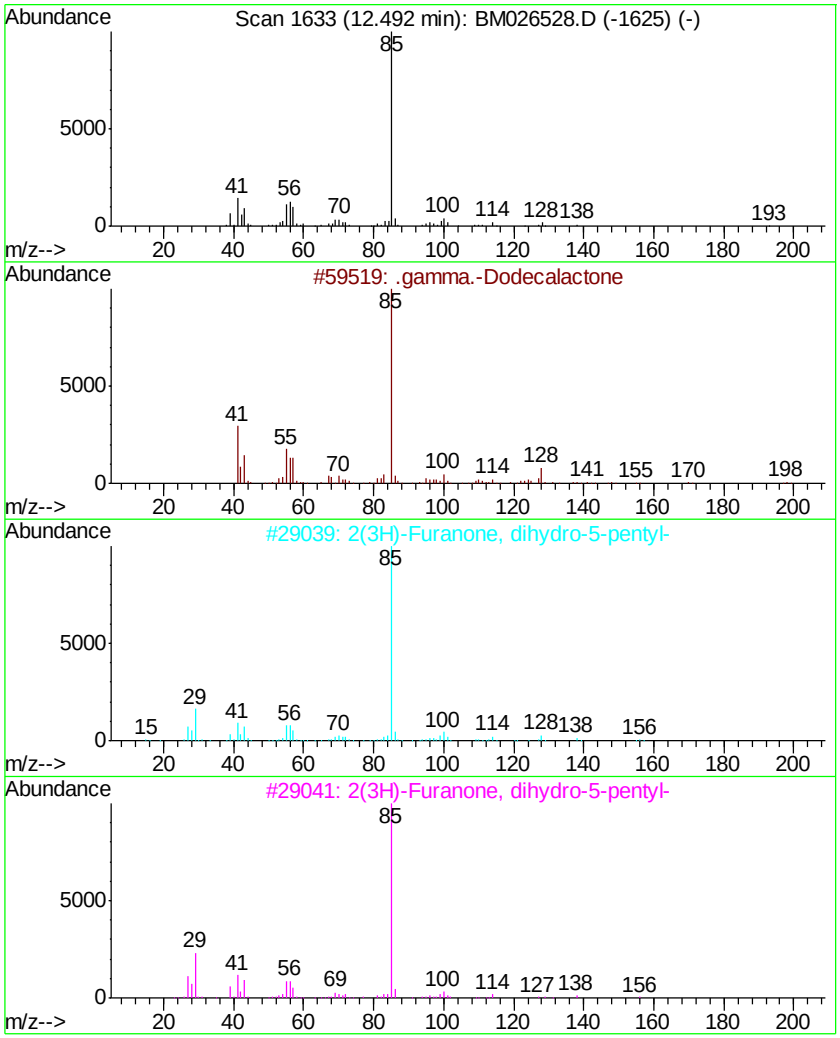
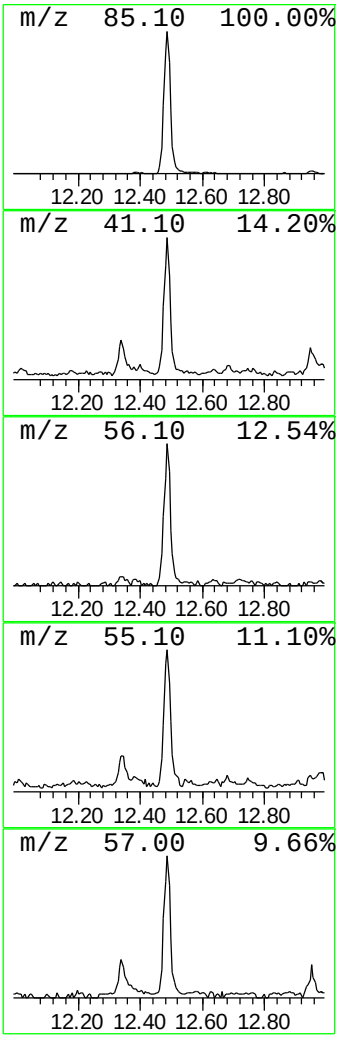
Peak Number 8 .gamma.-Dodecalactone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	2.74 ng/ul	235882	Acenaphthene-d10	14.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Dodecalactone	198	C12H22O2	002305-05-7	78
2	2(3H)-Furanone, dihydro-5-pentyl-	156	C9H16O2	000104-61-0	78
3	2(3H)-Furanone, dihydro-5-pentyl-	156	C9H16O2	000104-61-0	72
4	Pentanal, 5-[(tetrahydro-2H-pyra...	186	C10H18O3	014194-86-6	72
5	2(3H)-Furanone, 5-butyldihydro-	142	C8H14O2	000104-50-7	72





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Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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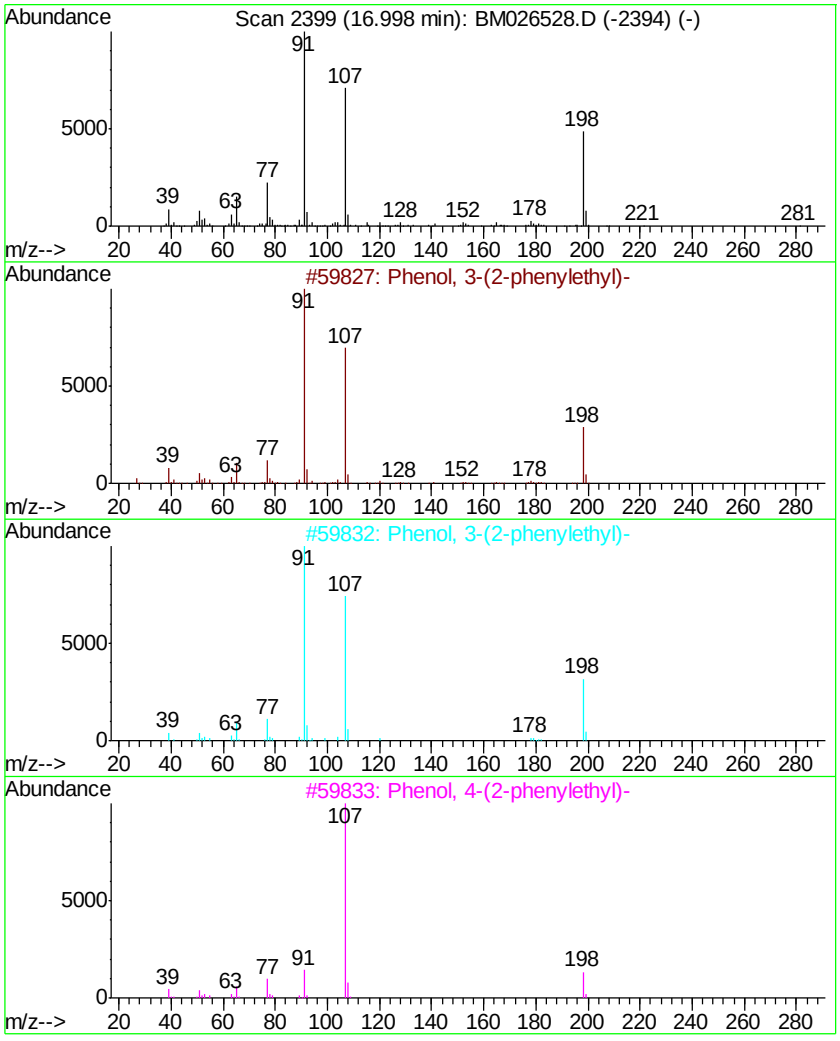
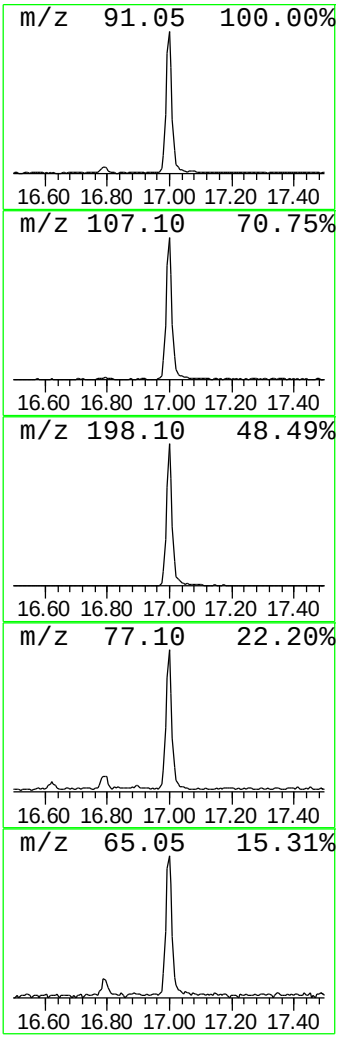
Peak Number 9 Phenol, 3-(2-phenylethyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	3.79 ng/ul	347652	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	93
2	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	90
3	Phenol, 4-(2-phenylethyl)-	198	C14H14O	006335-83-7	83
4	Propanedinitrile, (1-methylethen...	196	C13H12N2	069564-95-0	60
5	Benzyl mandelate	242	C15H14O3	000890-98-2	47





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

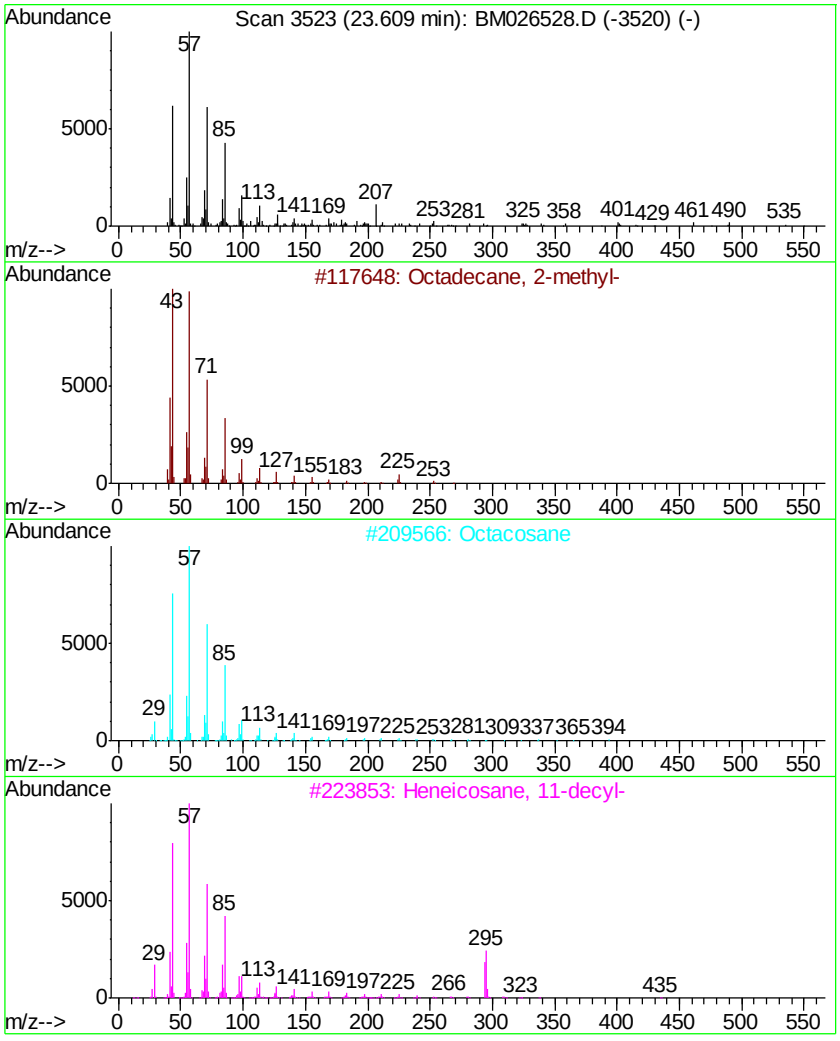
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

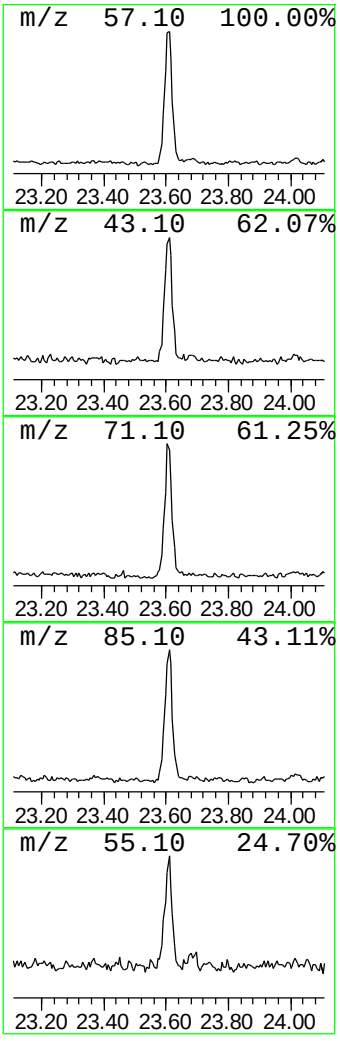
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	2.09 ng/ul	195418	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
2	Octacosane	394	C28H58	000630-02-4	81
3	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	81
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	81
5	Tetratetracontane	619	C44H90	007098-22-8	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062320\
Data File : BM026528.D
Acq On : 24 Jun 2020 10:09
Operator : JU/CG
Sample : L3069-10DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7G8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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
p-Cymene	7.52	9.1	ng/ul	302908	1	7.37	665753	20.0
Bicyclo[2.2.1]hep...	8.60	6.0	ng/ul	201195	1	7.37	665753	20.0
Cyclohexanemethan...	9.52	2.9	ng/ul	150534	2	10.13	1038700	20.0
Bicyclo[2.2.1]hep...	9.56	7.8	ng/ul	407071	2	10.13	1038700	20.0
Terpinen-4-ol	10.02	2.8	ng/ul	142928	2	10.13	1038700	20.0
.alpha.-Terpineol	10.21	4.7	ng/ul	244850	2	10.13	1038700	20.0
Hydrocinnamic acid	12.02	4.3	ng/ul	223227	2	10.13	1038700	20.0
.gamma.-Dodecalac...	12.49	2.7	ng/ul	235882	3	14.03	1724140	20.0
Phenol, 3-(2-phen...	17.00	3.8	ng/ul	347652	4	16.79	1832820	20.0
(DEL) Alkane: Str...	23.61	2.1	ng/ul	195418	6	23.09	1871690	20.0