LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

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-BM031420MA.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	4.181	216	220	225	rBV	38033	45411	0.19%	0.088%
2	4.240	226	230	242	rVB	38757	62946	0.26%	0.121%
3	4.816	324	328	341	rVB	183548	269409	1.12%	0.519%
4	5.275	400	406	413	rVB4	23799	40004	0.17%	0.077%
5	5.722	478	482	488	rBV3	15117	25765	0.11%	0.050%
6	6.422	596	601	607	rBV	57467	87484	0.36%	0.169%
7	6.934	684	688	695	rBV	37919	57400	0.24%	0.111%
8	7.198	727	733	738	rBV	748508	1077324	4.47%	2.077%
9	7.251	738	742	749	rVV	795596	1164788	4.83%	2.246%
10	7.316	749	753	761	rVB2	53891	95956	0.40%	0.185%
11	7.434	768	773	782	rBV	667981	1074450	4.46%	2.072%
12	7.639	801	808	816	rBV	697540	1131081	4.69%	2.181%
13	7.716	816	821	825	rVB	58056	85813	0.36%	0.165%
14	7.875	841	848	857	rBV	1027275	1549059	6.43%	2.987%
15	8.063	873	880	886	rBV	307375	473460	1.96%	0.913%
16	8.116	886	889	896	rVB3	20721	33111	0.14%	0.064%
17	8.257	908	913	922	rVB8	10951	26014	0.11%	0.050%
18	8.445	940	945	952	rVB2	60589	92404	0.38%	0.178%
19	8.769	992	1000	1006	rBV5	59558	132430	0.55%	0.255%
20	8.881	1013	1019	1024	rBV3	27516	50022	0.21%	0.096%
21	9.128	1057	1061	1066	rVB3	29309	42958	0.18%	0.083%
22	9.216	1071	1076	1080	rVB2	27805	39917	0.17%	0.077%
23	9.345	1094	1098	1108	rVB	73707	116064	0.48%	0.224%
24	9.433	1108	1113	1117	rBV	54224	76703	0.32%	0.148%
25	9.492	1117	1123	1132	rVB	76046	118778	0.49%	0.229%
26	9.580	1132	1138	1145	rVB	98440	150827	0.63%	0.291%
27	9.663	1145	1152	1156	rBV2	55426	101083	0.42%	0.195%
28	9.751	1163	1167	1174	rVB3	42528	78978	0.33%	0.152%
29	9.816	1174	1178	1182	rBV5	14091	27294	0.11%	0.053%
30	9.916	1188	1195	1200	rBV	87111	153667	0.64%	0.296%
31	10.045	1212	1217	1224	rVV9	13407	27094	0.11%	0.052%
32	10.122	1224	1230	1234	rVB5	25536	39931	0.17%	0.077%
33	10.292	1253	1259	1265	rVB4	26974	53547	0.22%	0.103%
34	10.404	1271	1278	1288	rVV	940501	1617610	6.71%	3.119%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

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-BM031420MA.M
Title : SVOA CALIBRATION

35	10.480	1288	1291	1297	rVB5	26895	41857	0.17%	0.081%
36	10.545	1297	1302	1307	rBV2	47524	77737	0.32%	0.150%
37	10.733	1332	1334	1340	rVB5	19220	27843	0.12%	0.054%
38	10.804	1340	1346	1354	rVB2	160161	253374	1.05%	0.489%
39	11.169	1400	1408	1413	rBV9	13778	29407	0.12%	0.057%
40	11.316	1427	1433	1437	rBV7	20022	38712	0.16%	0.075%
41	11.374	1437	1443	1452	rVB6	21248	52487	0.22%	0.101%
42	11.692	1493	1497	1502	rVB	97490	132365	0.55%	0.255%
43	11.833	1514	1521	1524	rBV3	26038	45132	0.19%	0.087%
44	11.892	1526	1531	1535	rVV2	84186	135162	0.56%	0.261%
45	11.963	1537	1543	1548	rVV3	45675	85229	0.35%	0.164%
46	12.039	1548	1556	1559	rVV5	40482	80125	0.33%	0.154%
47	12.074	1559	1562	1566	rVB3	21172	29998	0.12%	0.058%
48	12.180	1575	1580	1585	rBV	25154	42338	0.18%	0.082%
49	12.263	1587	1594	1598	rBV7	21226	51583	0.21%	0.099%
50	12.374	1609	1613	1619	rBV5	13649	27831	0.12%	0.054%
51	12.533	1634	1640	1641	rBV5	32233	54441	0.23%	0.105%
52	12.557	1642	1644	1647	rVB3	25685	27201	0.11%	0.052%
53	12.692	1662	1667	1673	rVB2	55084	85004	0.35%	0.164%
54	12.845	1688	1693	1697	rVB	40487	57759	0.24%	0.111%
55	12.945	1705	1710	1713	rVV3	29586	41955	0.17%	0.081%
56	13.027	1713	1724	1729	rVB3	82000	146375	0.61%	0.282%
57	13.221	1751	1757	1762	rVV4	21814	42315	0.18%	0.082%
58	13.280	1763	1767	1770	rVV5	22842	37673	0.16%	0.073%
59	13.327	1770	1775	1780	rVB2	41876	67580	0.28%	0.130%
60	13.463	1789	1798	1806	rBV7	25368	71946	0.30%	0.139%
61	13.545	1806	1812	1815	rBV7	26690	43993	0.18%	0.085%
62	13.627	1822	1826	1833	rBV2	58021	89916	0.37%	0.173%
63	13.686	1833	1836	1840	rVB	20139	24638	0.10%	0.048%
64	13.880	1863	1869	1876	rVV	16425	42081	0.17%	0.081%
65	13.957	1876	1882	1886	rVV3	25644	52425	0.22%	0.101%
66	14.010	1886	1891	1897	rVB4	17090	34751	0.14%	0.067%
67	14.086	1900	1904	1909	rVV4	20709	27354	0.11%	0.053%
68	14.268	1928	1935	1942	rVB2	1521527	2201144	9.13%	4.244%
69	14.357	1946	1950	1953	rVB3	22009	29447	0.12%	0.057%
70	14.468	1964	1969	1976	rBV8	21150	38573	0.16%	0.074%
71	14.592	1982	1990	1995	rBV2	182976	367693	1.53%	0.709%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

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-BM031420MA.M
Title : SVOA CALIBRATION

72	14.639	1995	1998	2007	rVV6	63300	148694	0.62%	0.287%
73	14.715	2007	2011	2014	rVV	78834	121459	0.50%	0.234%
74	14.751	2014	2017	2023	rVV2	53205	77730	0.32%	0.150%
75	14.815	2024	2028	2033	rVV3	50696	80970	0.34%	0.156%
76	14.898	2033	2042	2055	rVB	1527924	2159666	8.96%	4.164%
77	15.086	2067	2074	2078	rVV	34947	58693	0.24%	0.113%
78	15.262	2100	2104	2109	rVB2	25477	38874	0.16%	0.075%
79	15.380	2120	2124	2127	rBV	29365	39790	0.17%	0.077%
80	15.921	2213	2216	2227	rBV10	8931	26917	0.11%	0.052%
81	16.098	2240	2246	2253	rBV	425268	555606	2.30%	1.071%
82	16.680	2337	2345	2362	rBV	17184908	24109041	100.00%	46.484%
83	17.015	2394	2402	2408	rBV	1927148	2704304	11.22%	5.214%
84	17.074	2408	2412	2415	rVV3	40832	59421	0.25%	0.115%
85	17.109	2415	2418	2422	rVB	40022	55232	0.23%	0.106%
86	17.404	2465	2468	2474	rVB8	26350	43106	0.18%	0.083%
87	19.409	2804	2809	2816	rBV	34032	55500	0.23%	0.107%
88	20.709	3027	3030	3035	rBV3	28196	39464	0.16%	0.076%
89	20.903	3060	3063	3067	rBV3	31485	36615	0.15%	0.071%
90	20.991	3076	3078	3082	rBV5	24051	27821	0.12%	0.054%
91	21.203	3108	3114	3119	rBV	2552593	3175718	13.17%	6.123%
92	23.380	3478	3484	3494	rVB	1896183	3296154	13.67%	6.355%

Sum of corrected areas: 51865001

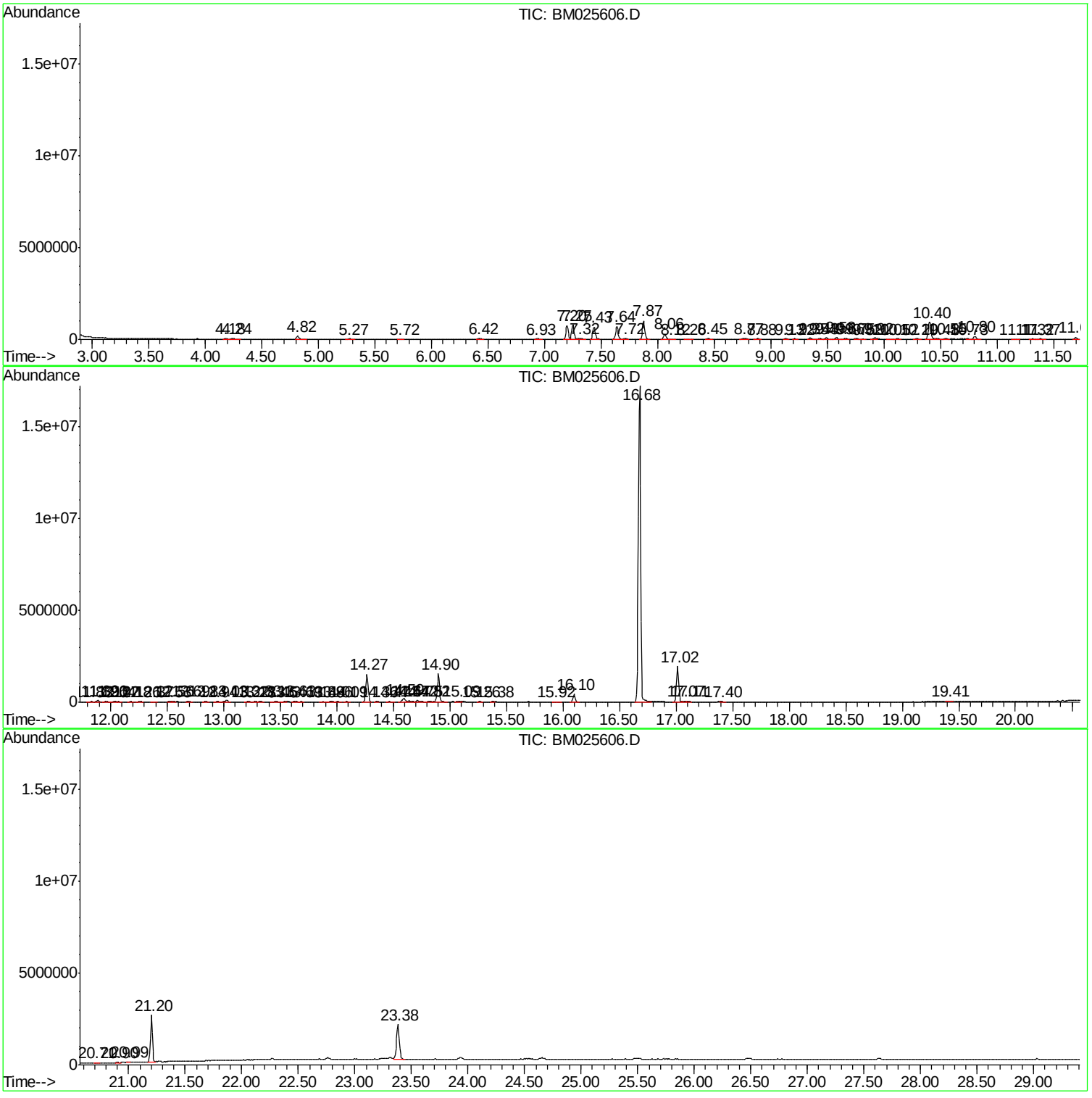
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.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\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

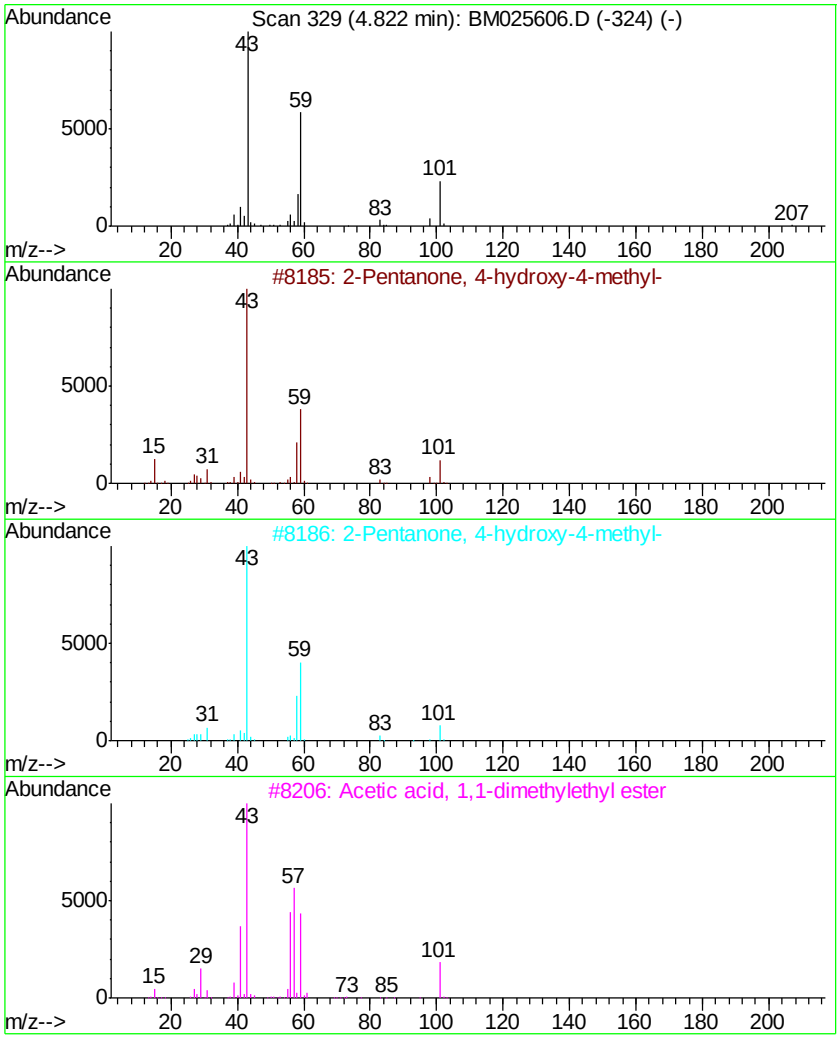
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

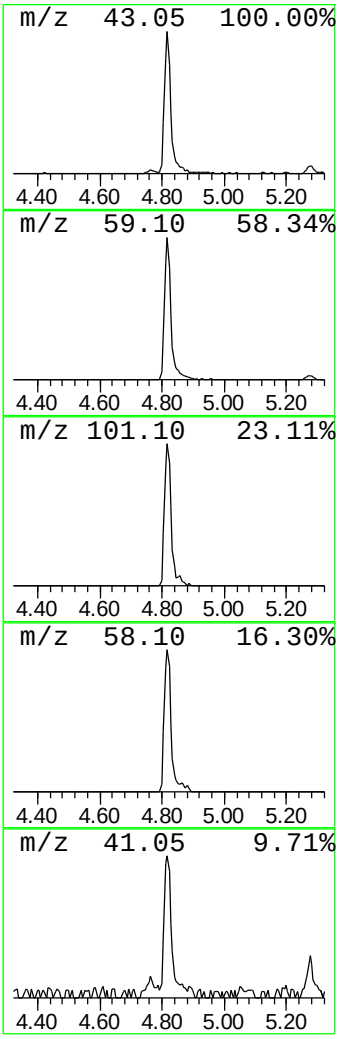
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	4.76 ng/ul	269409	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	32





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

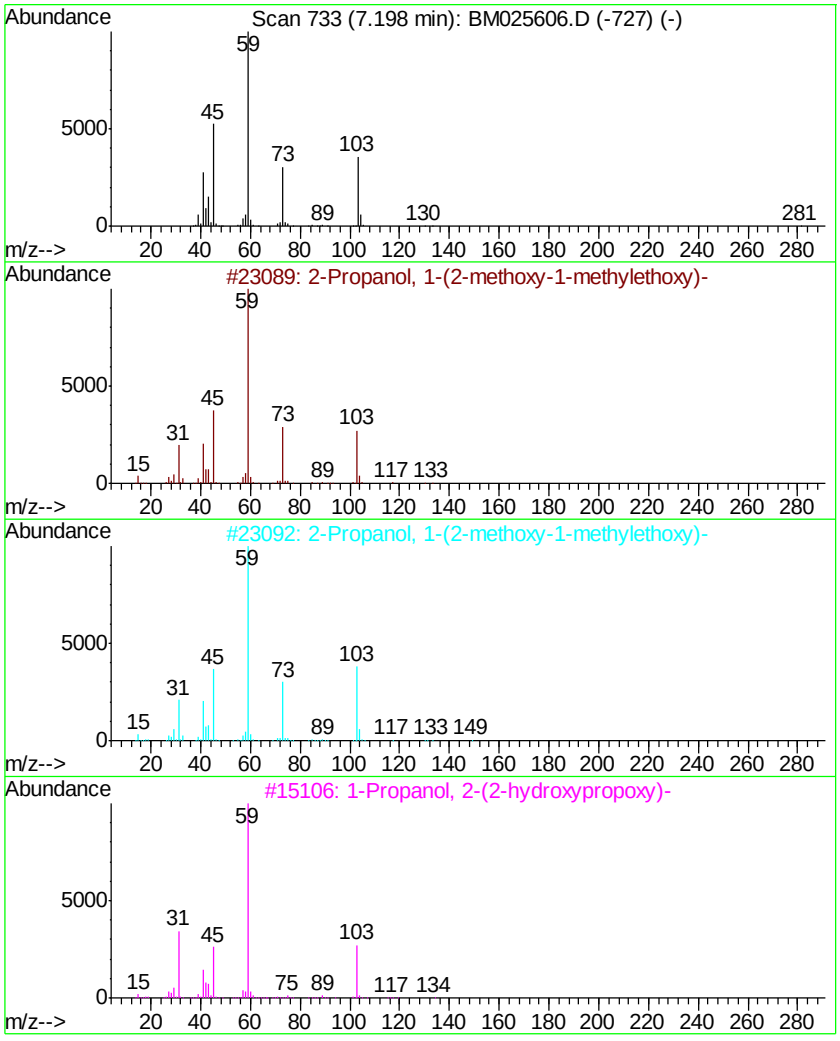
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

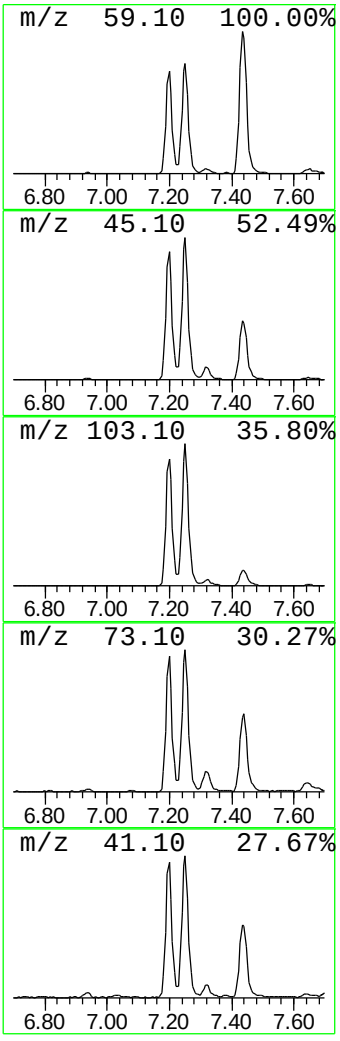
****

Peak Number 2 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	19.05 ng/ul	1077320	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	90
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
4	2-Propanol, 1-(2-ethoxypropoxy)-	162	C8H18O3	010143-32-5	56
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

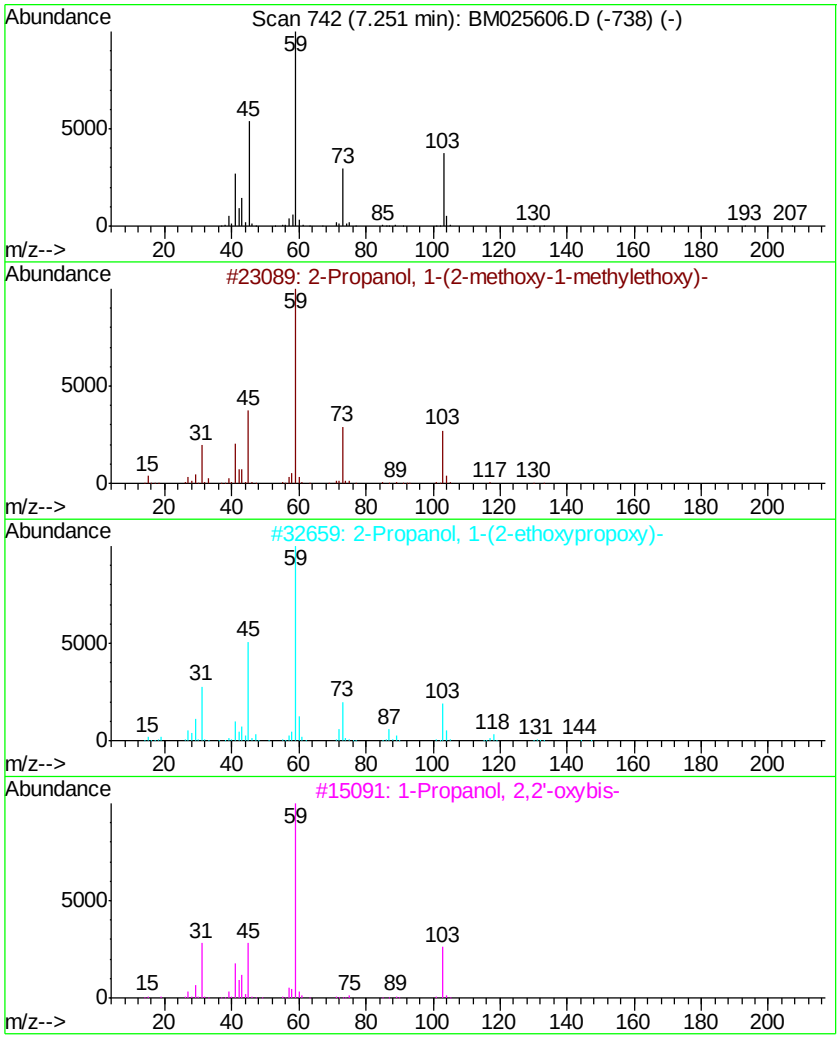
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

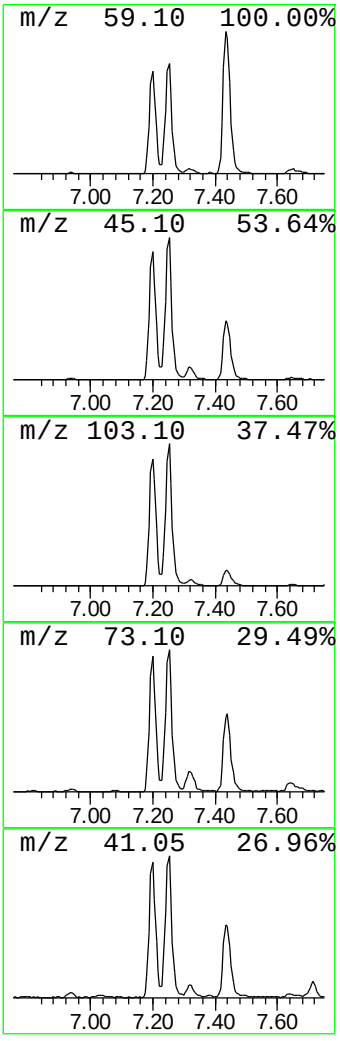
****

Peak Number 3 2-Propanol, 1-(2-ethoxyprop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	20.60 ng/ul	1164790	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
2	2-Propanol, 1-(2-ethoxypropoxy)-	162	C8H18O3	010143-32-5	56
3	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	53
4	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	45





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

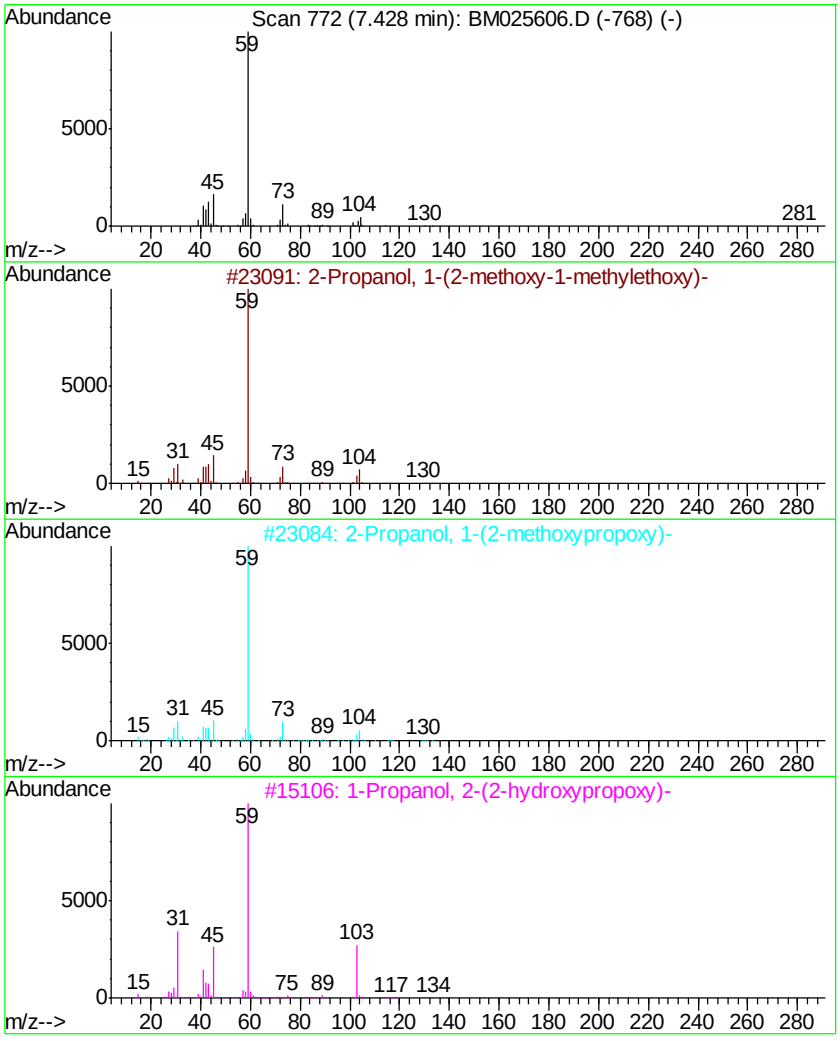
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

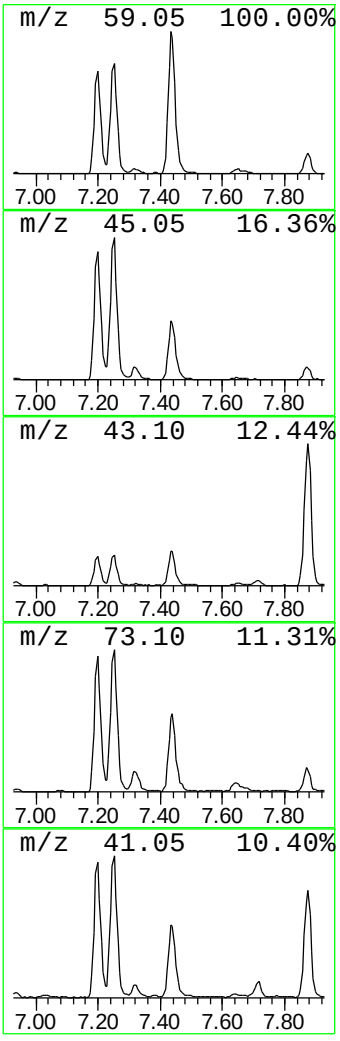
****

Peak Number 4 2-Propanol, 1-(2-methoxypro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.43	19.00 ng/ul	1074450	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	83
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

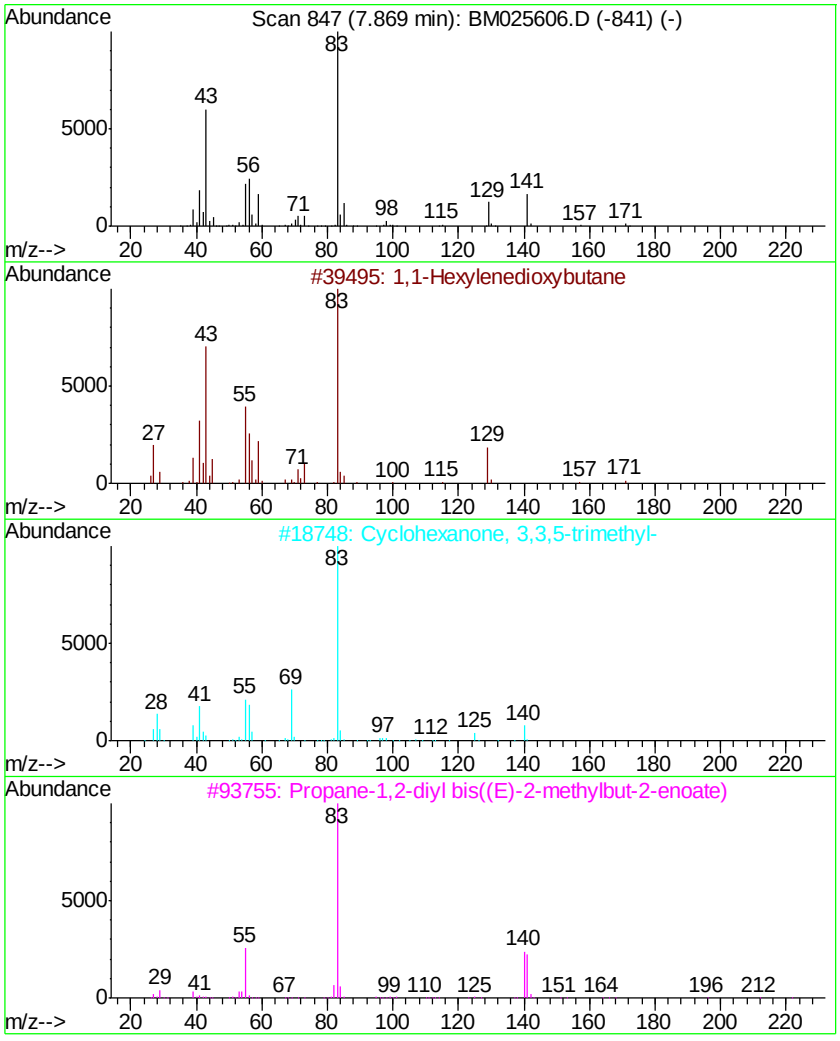
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

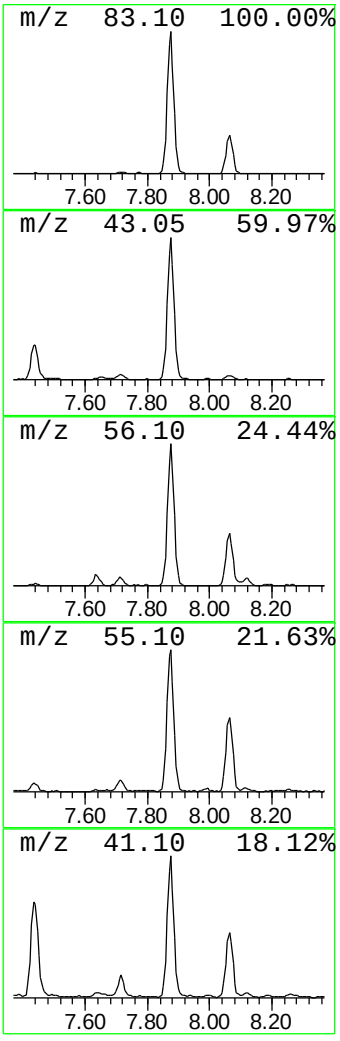
****

Peak Number 5 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	27.39 ng/ul	1549060	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	36
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	33
3	Propane-1,2-diyl bis((E)-2-methy...	240	C13H20O4	1000373-72-4	33
4	3-(Butylamino)propionitrile	126	C7H14N2	000693-51-6	28
5	Oxalic acid, cyclohexyl pentyl e...	242	C13H22O4	1000309-30-6	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

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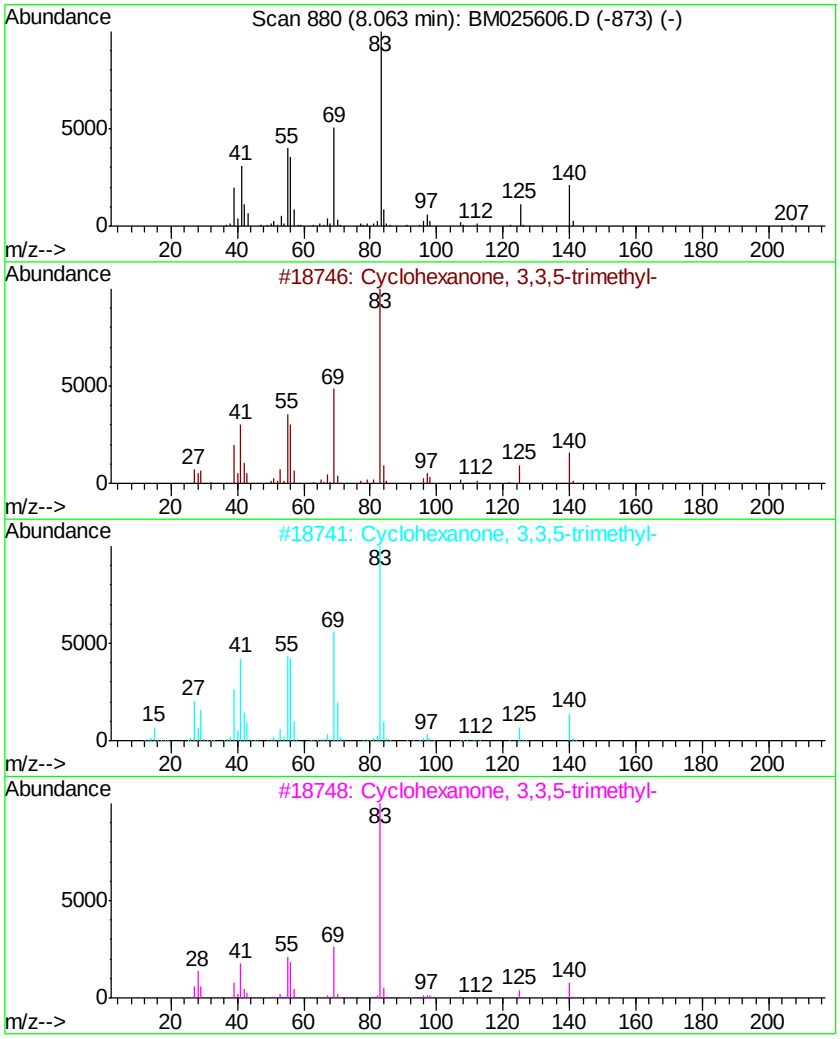
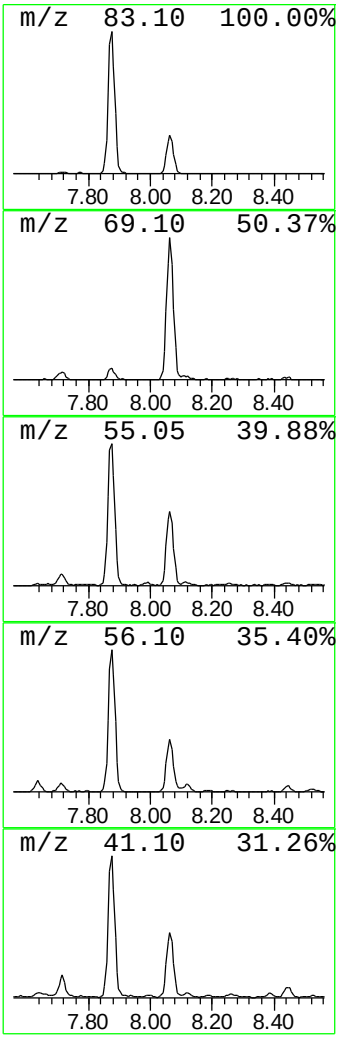
Peak Number 6 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	8.37 ng/ul	473460	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
4	Imidazole, 2-acetamido-	125	C5H7N3O	052737-49-2	58
5	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	52





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

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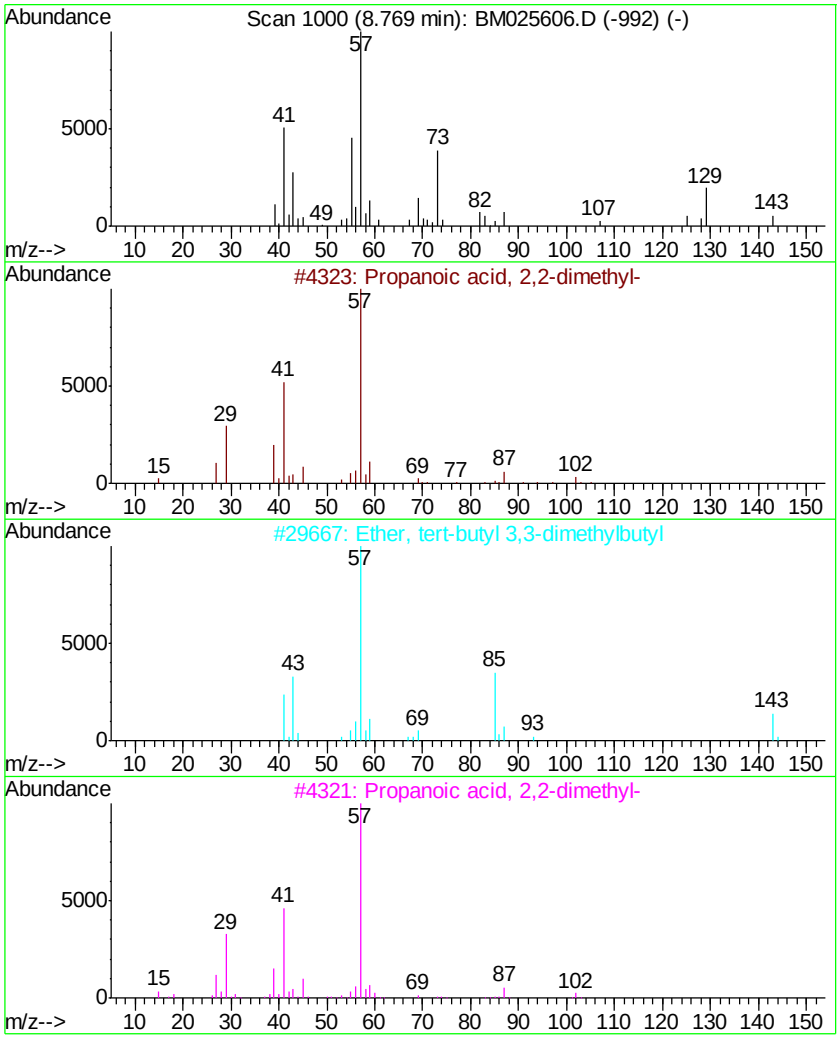
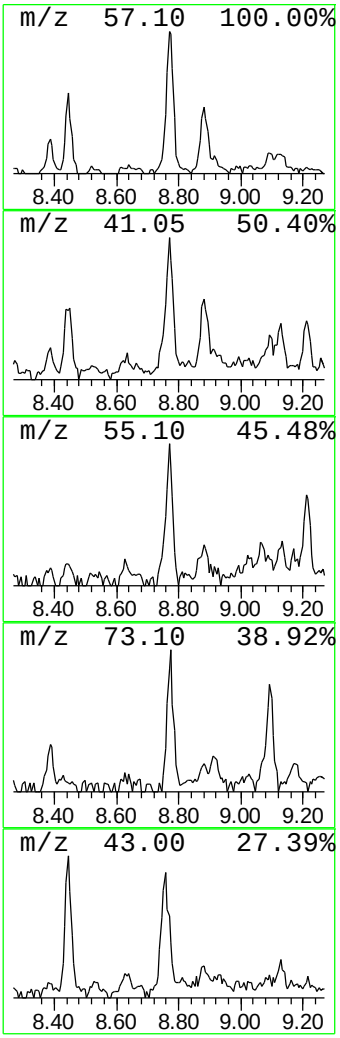
Peak Number 7 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.77	2.34 ng/ul	132430	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	35
2	Ether, tert-butyl 3,3-dimethylbutyl	158	C10H22O	004419-58-3	23
3	Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	17
4	Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	17
5	Ethynyl tert-butyl sulfoxide	130	C6H10OS	041463-34-7	16





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Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

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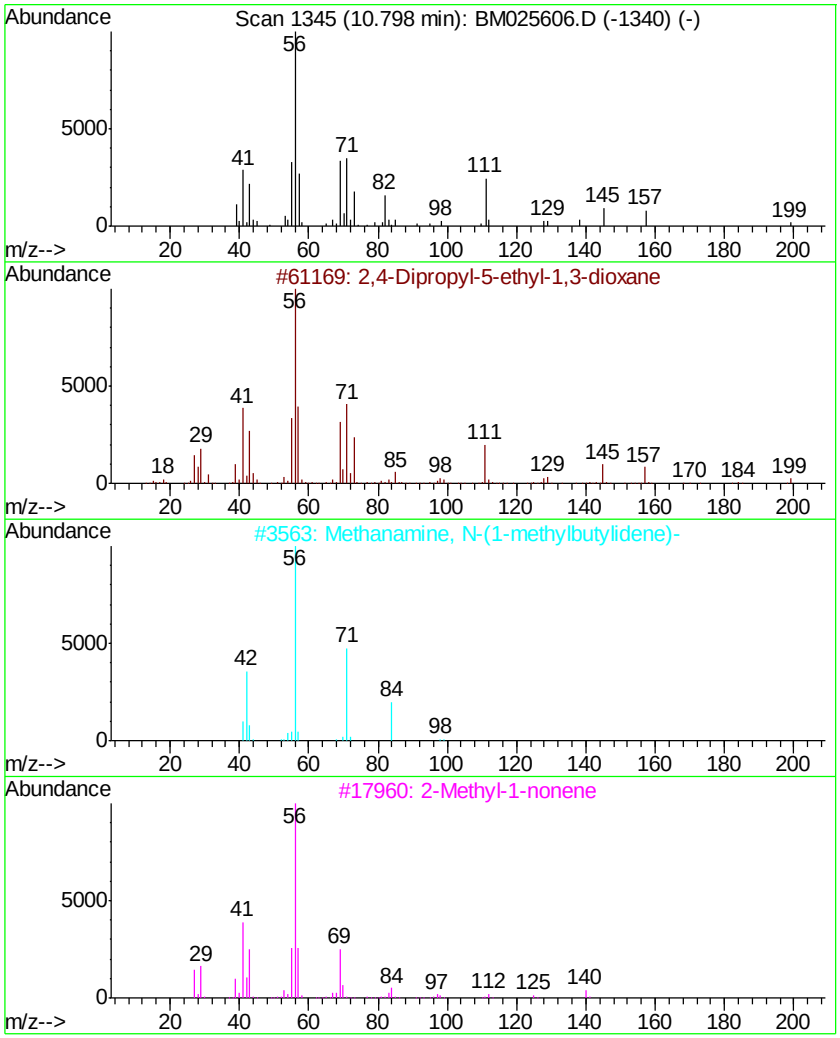
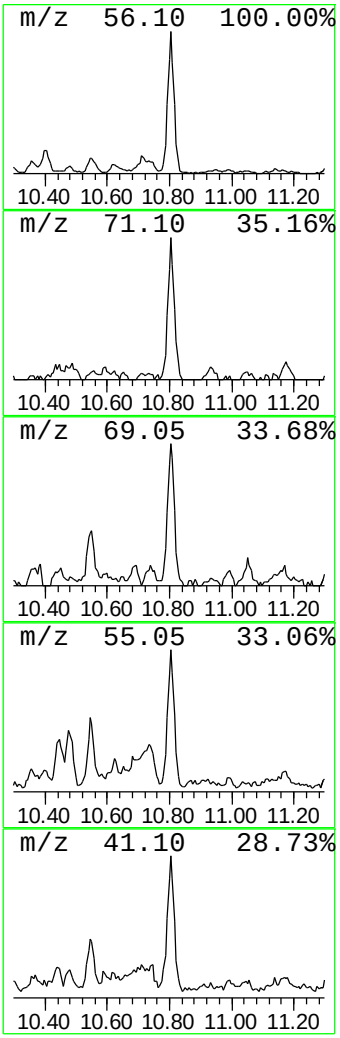
Peak Number 8 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	3.13 ng/ul	253374	Naphthalene-d8	10.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	50
2	Methanamine, N-(1-methylbutylide...	99	C6H13N	022431-09-0	38
3	2-Methyl-1-nonene	140	C10H20	002980-71-4	35
4	1-Octene, 2-methyl-	126	C9H18	004588-18-5	32
5	1-Octene, 2-methyl-	126	C9H18	004588-18-5	32





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

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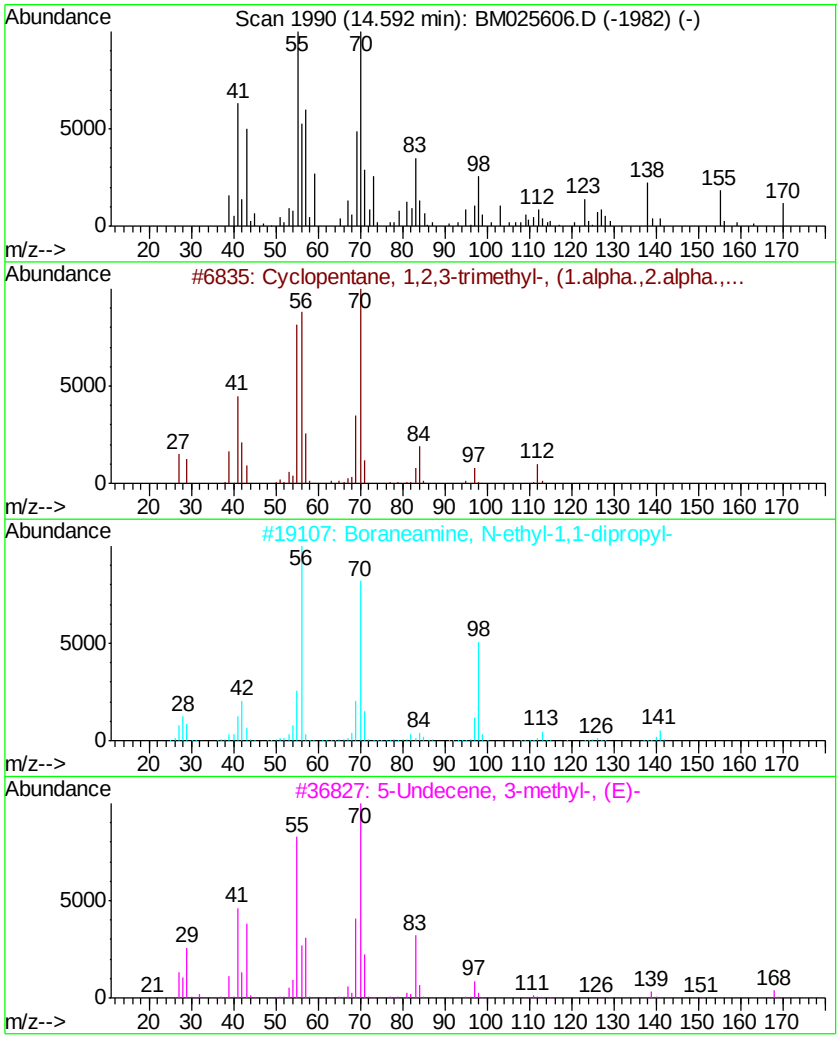
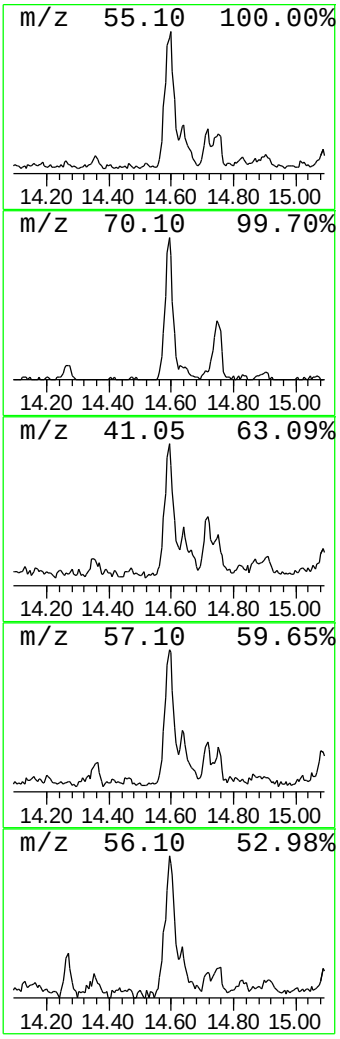
Peak Number 9 (DEL) Alkane: Cyclic14.59 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	3.34 ng/ul	367693	Acenaphthene-d10	14.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	015890-40-1	38
2	Boraneamine, N-ethyl-1,1-dipropyl-	141	C8H20BN	1000160-33-0	30
3	5-Undecene, 3-methyl-, (E)-	168	C12H24	074630-67-4	27
4	2-Heptene, 3-methyl-	112	C8H16	003404-75-9	22
5	3-Heptene	98	C7H14	000592-78-9	20





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

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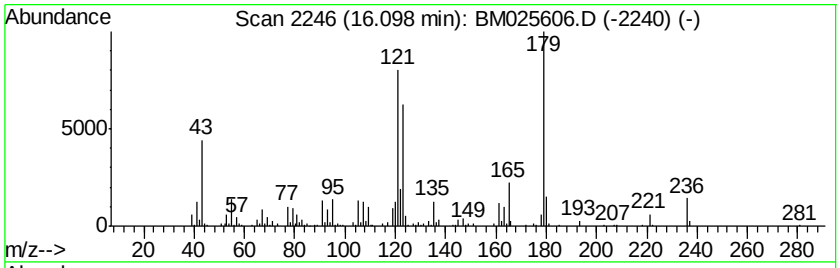
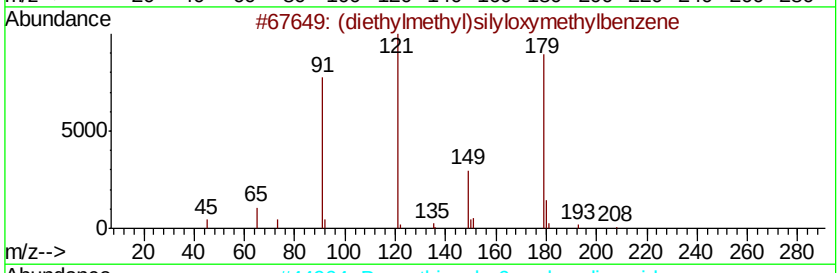
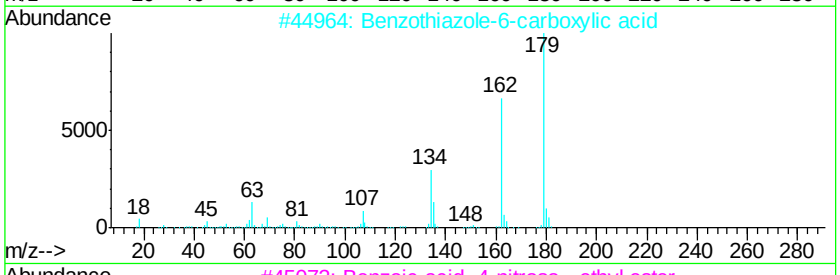
Peak Number 10 unknown-03 Concentration Rank 7

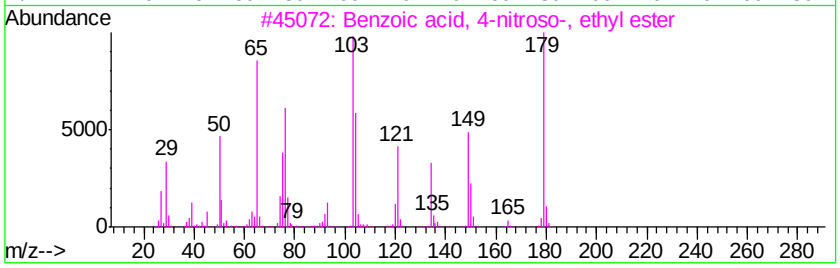
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	4.11 ng/ul	555606	Phenanthrene-d10	17.02

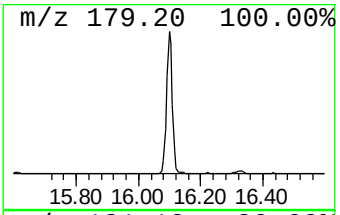
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(diethylmethyl)silyloxymethylben...	208	C12H20OSi	052313-30-1	47
2	Benzothiazole-6-carboxylic acid	179	C8H5NO2S	003622-35-3	30
3	Benzoic acid, 4-nitroso-, ethyl ...	179	C9H9NO3	007476-79-1	30
4	1-(4-Methoxybenzyloxy)-1-methyl-...	236	C13H20O2Si	1000280-11-2	27
5	(p-(Allyloxycarbonyl)phenoxy)ace...	236	C12H12O5	1000241-83-7	27

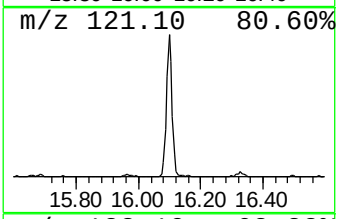


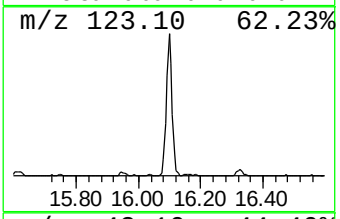


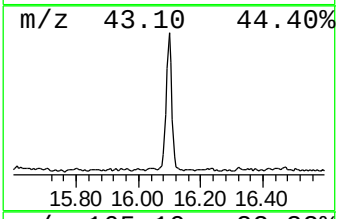


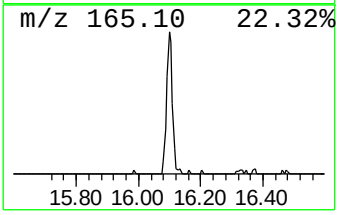












Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031620\
Data File : BM025606.D
Acq On : 17 Mar 2020 09:39
Operator : CG/JU
Sample : L1840-17DL 100X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBET2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.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...	4.82	4.8	ng/ul	269409	1	7.64	1131080	20.0
2-Propanol, 1-(2-...	7.20	19.1	ng/ul	1077320	1	7.64	1131080	20.0
2-Propanol, 1-(2-...	7.25	20.6	ng/ul	1164790	1	7.64	1131080	20.0
2-Propanol, 1-(2-...	7.43	19.0	ng/ul	1074450	1	7.64	1131080	20.0
unknown-01	7.87	27.4	ng/ul	1549060	1	7.64	1131080	20.0
Cyclohexanone, 3,...	8.06	8.4	ng/ul	473460	1	7.64	1131080	20.0
unknown-02	8.77	2.3	ng/ul	132430	1	7.64	1131080	20.0
2,4-Dipropyl-5-et...	10.80	3.1	ng/ul	253374	2	10.40	1617610	20.0
(DEL) Alkane: Cyc...	14.59	3.3	ng/ul	367693	3	14.27	2201140	20.0
unknown-03	16.10	4.1	ng/ul	555606	4	17.02	2704300	20.0