LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.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	3.175	10	15	19	rBV	92657	123251	3.66%	0.354%
2	3.287	29	34	38	rBV3	27010	35717	1.06%	0.103%
3	3.775	113	117	127	rBV2	60233	99275	2.95%	0.285%
4	3.910	134	140	142	rBV5	35137	51486	1.53%	0.148%
5	4.152	176	181	186	rBV	113046	154523	4.59%	0.444%
6	4.205	186	190	199	rVB	195974	271036	8.06%	0.778%
7	4.922	308	312	320	rVB5	23810	48117	1.43%	0.138%
8	5.316	371	379	387	rBV2	79081	146153	4.35%	0.419%
9	5.593	420	426	433	rBV	144763	215895	6.42%	0.620%
10	5.751	447	453	463	rVB4	39511	75759	2.25%	0.217%
11	5.887	472	476	482	rVB	37039	55760	1.66%	0.160%
12	6.187	521	527	530	rBV5	22027	39411	1.17%	0.113%
13	6.304	541	547	552	rBV4	26205	45992	1.37%	0.132%
14	6.463	568	574	576	rBV4	30838	45892	1.36%	0.132%
15	6.851	633	640	644	rBV	85750	141070	4.19%	0.405%
16	6.987	654	663	671	rVB2	120053	219920	6.54%	0.631%
17	7.140	683	689	694	rBV	402579	616783	18.34%	1.770%
18	7.175	694	695	703	rVV2	36384	47553	1.41%	0.136%
19	7.245	703	707	714	rVB2	34472	63138	1.88%	0.181%
20	7.340	717	723	732	rBV	431488	700296	20.82%	2.010%
21	7.804	795	802	807	rBV	468911	782531	23.27%	2.246%
22	8.351	892	895	900	rBV2	20349	36510	1.09%	0.105%
23	8.410	900	905	915	rVB3	89879	181611	5.40%	0.521%
24	8.522	918	924	930	rBV2	204598	337228	10.03%	0.968%
25	8.634	939	943	950	rVB2	49197	80446	2.39%	0.231%
26	8.975	994	1001	1007	rBV	555629	953677	28.35%	2.737%
27	9.034	1008	1011	1017	rVV5	34801	61970	1.84%	0.178%
28	9.098	1017	1022	1028	rVB2	53493	92621	2.75%	0.266%
29	9.163	1028	1033	1037	rBV4	21537	33696	1.00%	0.097%
30	9.516	1088	1093	1100	rBV7	19095	35630	1.06%	0.102%
31	9.692	1116	1123	1134	rBV	462999	793940	23.60%	2.279%
32	9.798	1134	1141	1150	rVB	35089	69359	2.06%	0.199%
33	9.939	1153	1165	1173	rBV2	121770	239258	7.11%	0.687%
34	10.234	1208	1215	1224	rBV2	536624	946961	28.15%	2.718%

LSC Area Percent Report

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Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Title : SVOA CALIBRATION

35	10.604	1265	1278	1290	rBV	604992	1252453	37.24%	3.595%
36	10.763	1299	1305	1307	rBV5	21757	39905	1.19%	0.115%
37	10.992	1338	1344	1349	rBV8	16249	38937	1.16%	0.112%
38	11.404	1408	1414	1428	rBV5	81103	185141	5.50%	0.531%
39	11.510	1428	1432	1436	rVV4	17044	34494	1.03%	0.099%
40	11.563	1436	1441	1446	rVB6	30216	50592	1.50%	0.145%
41	11.939	1499	1505	1512	rBV8	34954	73298	2.18%	0.210%
42	12.069	1522	1527	1531	rBV5	39528	67450	2.01%	0.194%
43	12.222	1547	1553	1556	rVV6	23973	43409	1.29%	0.125%
44	12.310	1560	1568	1577	rBV3	39791	93371	2.78%	0.268%
45	12.727	1634	1639	1645	rBV6	26787	53927	1.60%	0.155%
46	12.792	1645	1650	1656	rVB6	24856	47475	1.41%	0.136%
47	12.904	1665	1669	1677	rVB7	18434	34353	1.02%	0.099%
48	13.039	1689	1692	1696	rBV3	33243	45931	1.37%	0.132%
49	13.098	1697	1702	1706	rVB2	53561	78681	2.34%	0.226%
50	13.351	1738	1745	1752	rBV3	52769	91156	2.71%	0.262%
51	13.863	1826	1832	1837	rBV	1245507	1697892	50.48%	4.873%
52	14.151	1875	1881	1887	rVB	1251707	1836045	54.59%	5.270%
53	14.457	1927	1933	1942	rVB2	931825	1433283	42.61%	4.114%
54	14.704	1966	1975	1983	rBV4	101693	208075	6.19%	0.597%
55	15.274	2067	2072	2079	rVB	89418	132720	3.95%	0.381%
56	15.451	2096	2102	2110	rBV	1675560	2358886	70.13%	6.770%
57	15.592	2120	2126	2135	rBV	582853	874719	26.01%	2.511%
58	16.739	2316	2321	2330	rVB4	45084	74460	2.21%	0.214%
59	17.010	2362	2367	2372	rBV2	30438	44422	1.32%	0.127%
60	17.210	2395	2401	2408	rVB	1326379	1841681	54.75%	5.286%
61	17.310	2411	2418	2425	rBV	1834592	2587564	76.93%	7.427%
62	18.157	2560	2562	2570	rBV2	28440	41361	1.23%	0.119%
63	18.857	2676	2681	2688	rBV	249334	331106	9.84%	0.950%
64	19.239	2743	2746	2749	rVB4	30901	35860	1.07%	0.103%
65	19.492	2786	2789	2793	rBV6	25566	39986	1.19%	0.115%
66	19.604	2803	2808	2813	rBV	2436197	3363549	100.00%	9.654%
67	19.645	2813	2815	2821	rVB	146659	211832	6.30%	0.608%
68	20.515	2961	2963	2967	rBV4	54532	60734	1.81%	0.174%
69	21.392	3108	3112	3117	rBV	1986216	2429852	72.24%	6.974%
70	23.556	3475	3480	3489	rVB2	1605229	3009427	89.47%	8.637%
71	23.703	3500	3505	3515	rVB	1124072	2155013	64.07%	6.185%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Title : SVOA CALIBRATION

Sum of corrected areas: 34841505

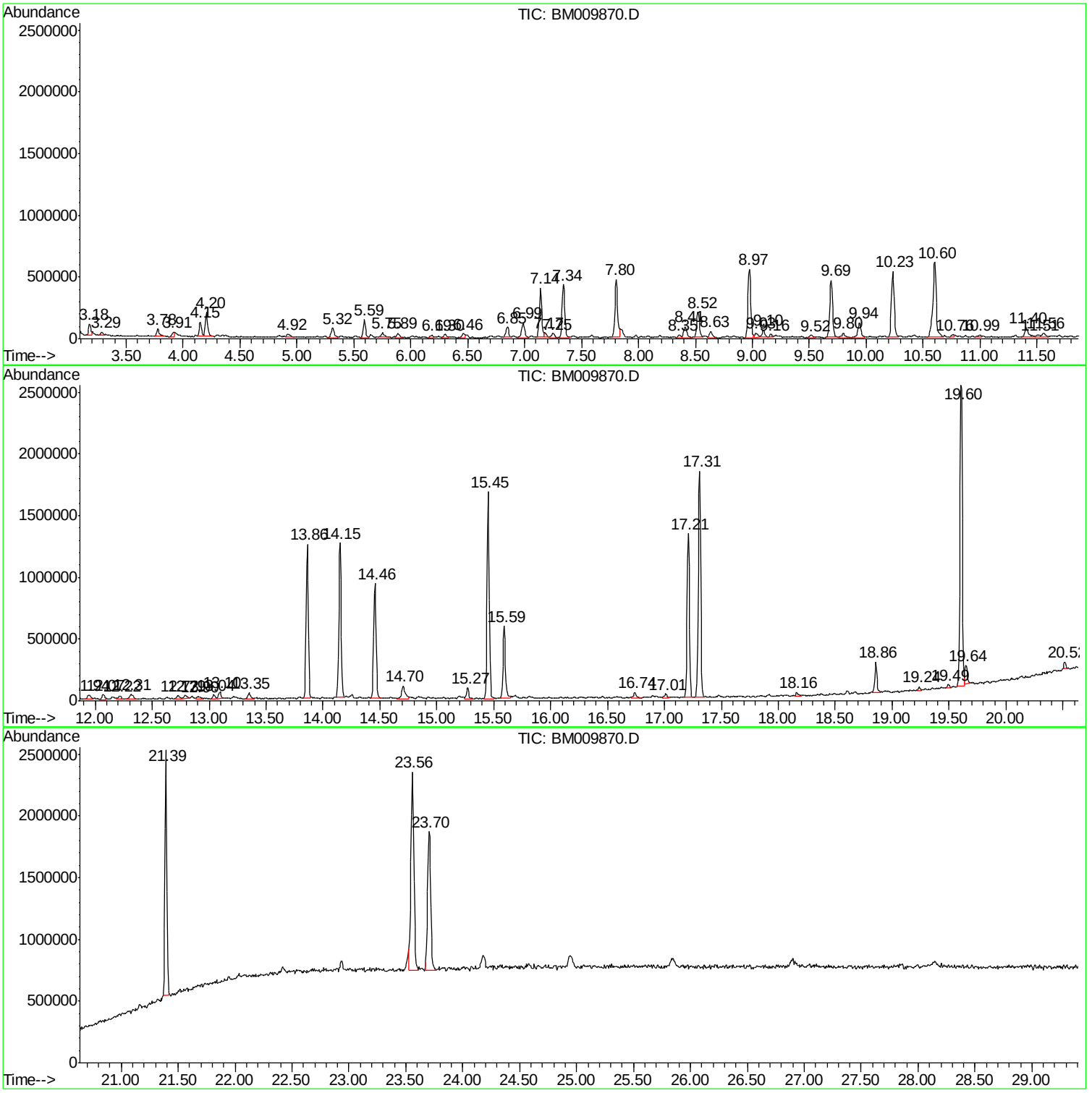
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

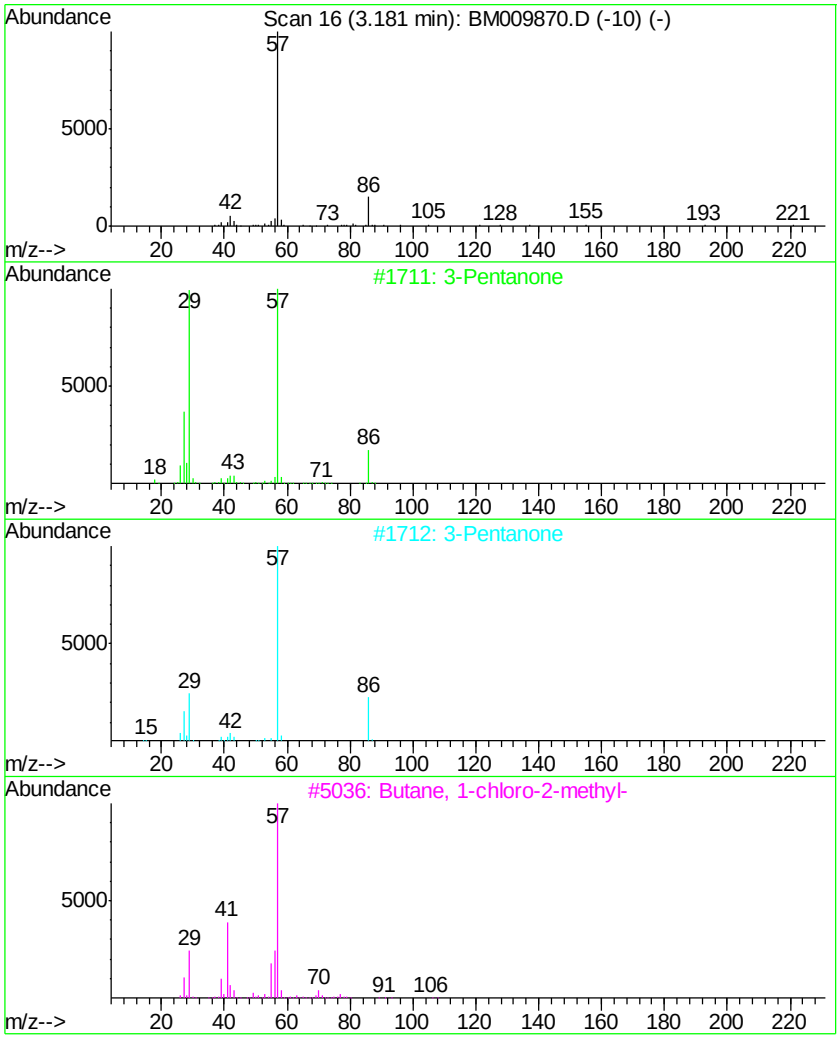
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

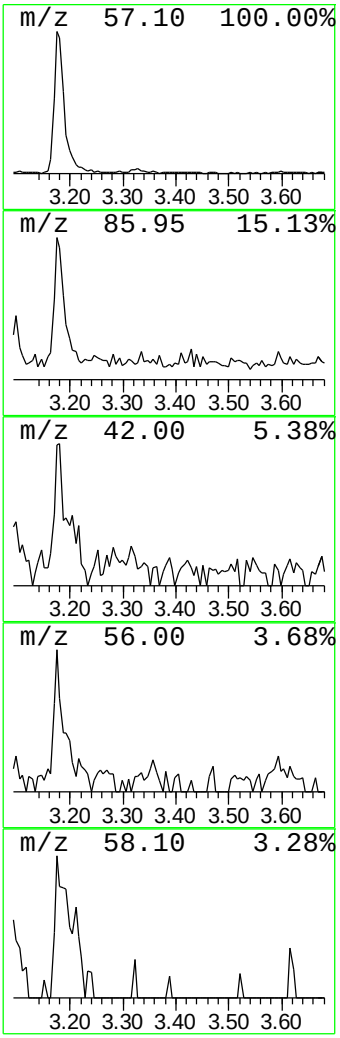
****

Peak Number 1 3-Pentanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	3.15 ng/ul	123251	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone	86	C5H10O	000096-22-0	58
2	3-Pentanone	86	C5H10O	000096-22-0	53
3	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	12
4	3-Pentanone	86	C5H10O	000096-22-0	9
5	Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	9





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

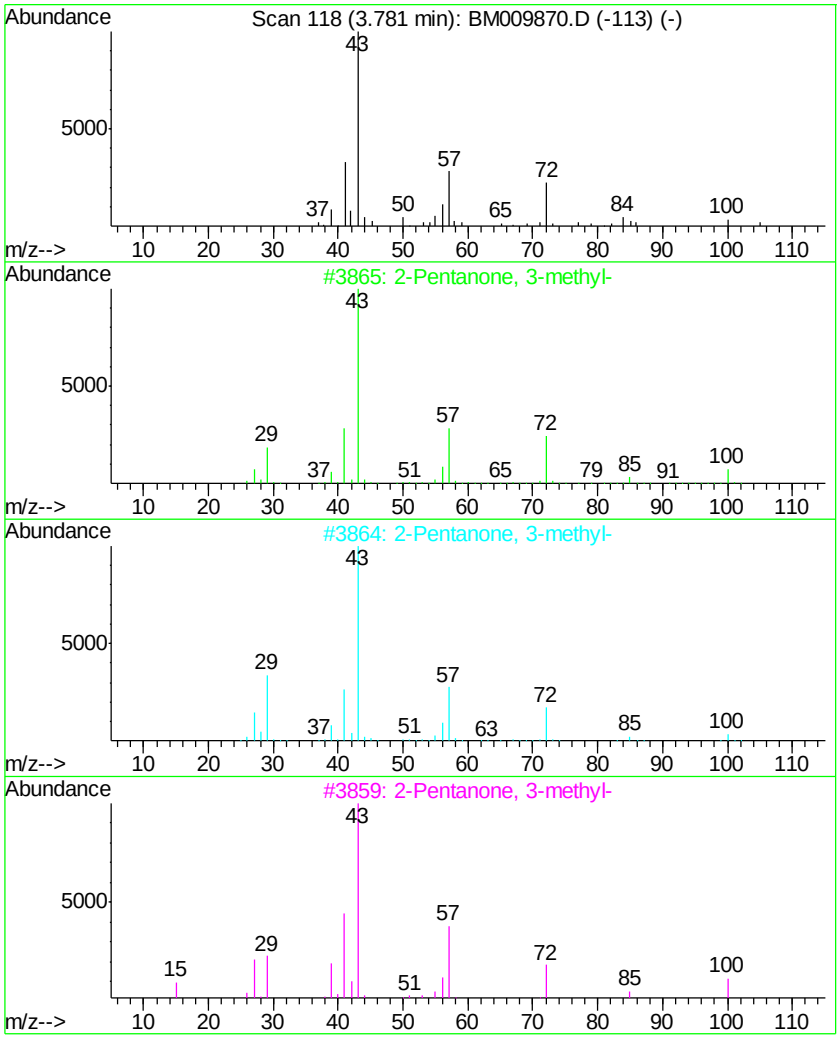
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

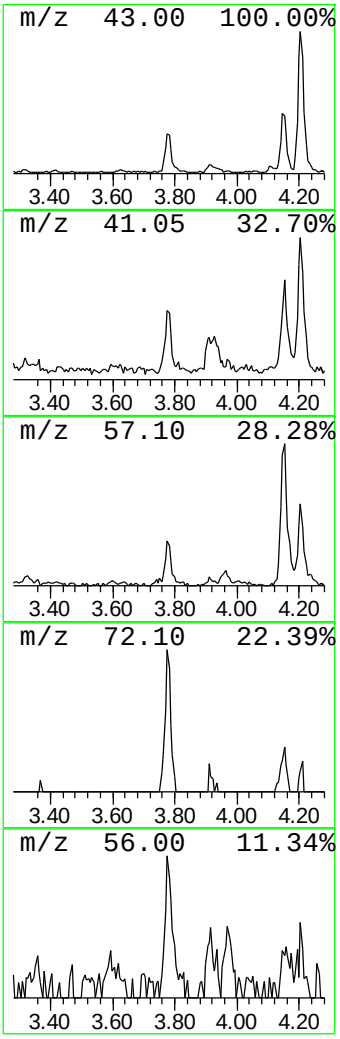
****

Peak Number 2 2-Pentanone, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.78	2.54 ng/ul	99275	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	74
2	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	72
3	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	50
4	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	50
5	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	50





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

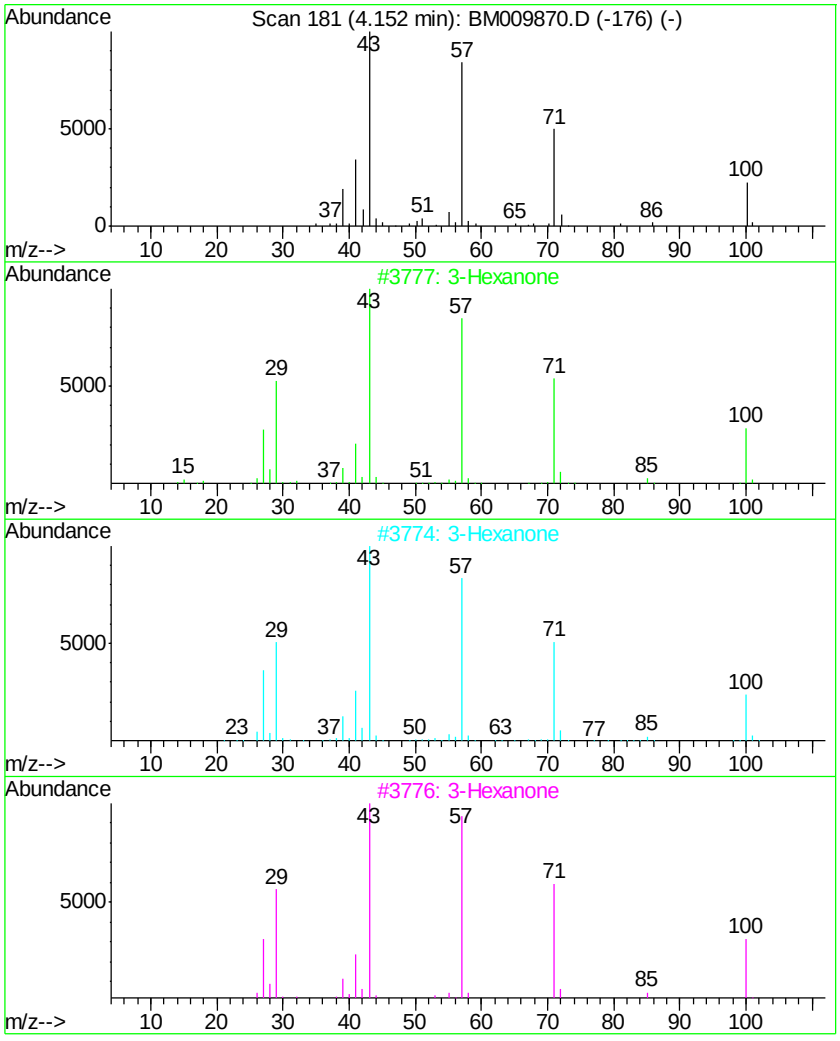
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

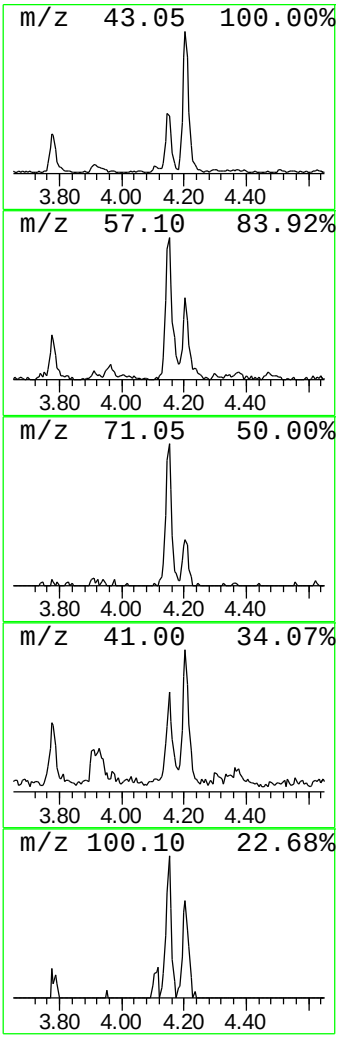
****

Peak Number 3 3-Hexanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.15	3.95 ng/ul	154523	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	86
2	3-Hexanone	100	C6H12O	000589-38-8	86
3	3-Hexanone	100	C6H12O	000589-38-8	72
4	3-Hexanone	100	C6H12O	000589-38-8	59
5	3-Hexanone	100	C6H12O	000589-38-8	43





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

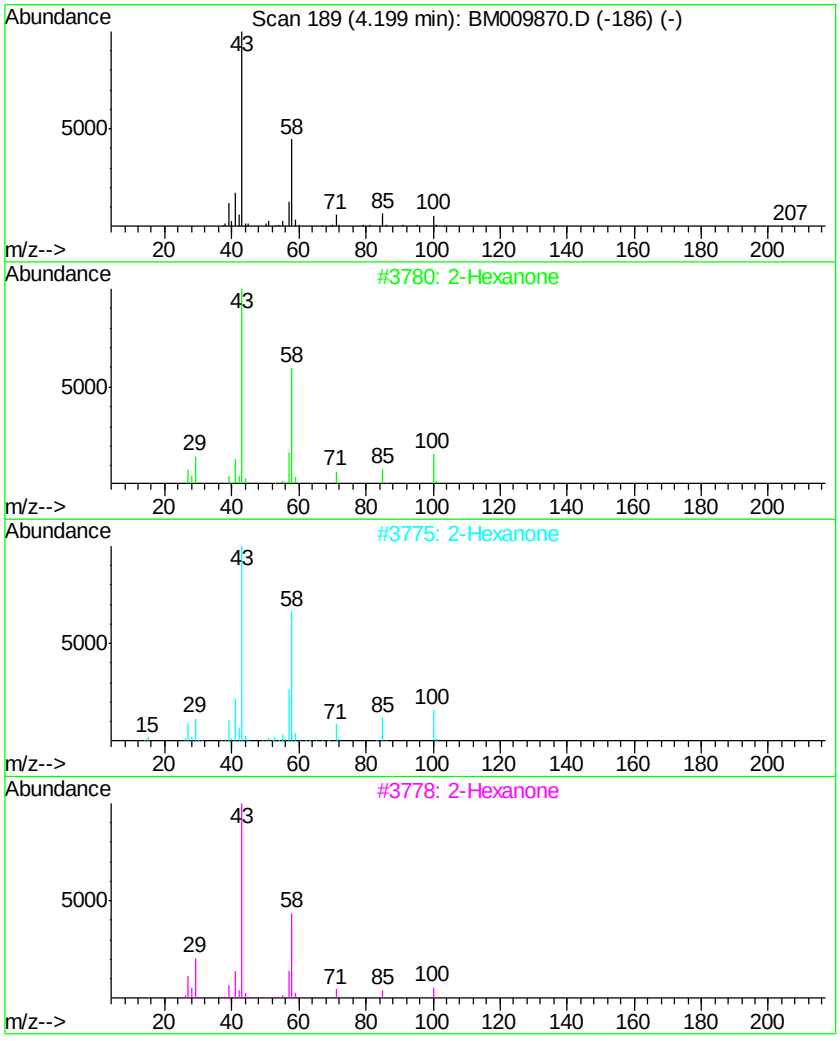
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

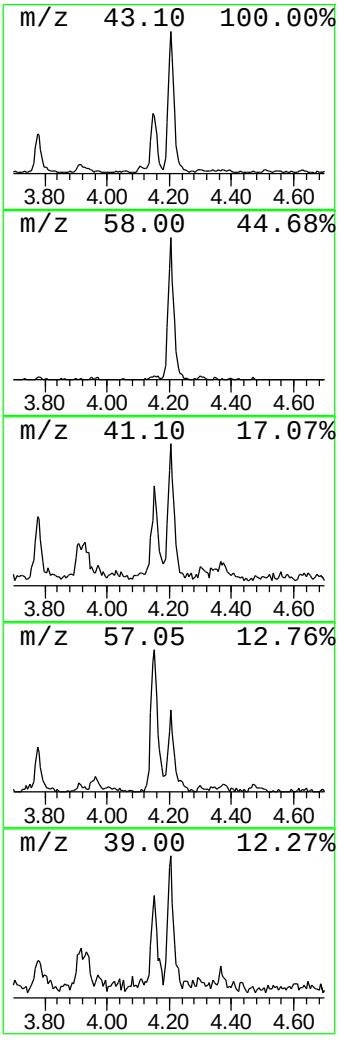
****

Peak Number 4 2-Hexanone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	6.93 ng/ul	271036	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	2-Hexanone	100	C6H12O	000591-78-6	90
3	2-Hexanone	100	C6H12O	000591-78-6	86
4	2-Hexanone	100	C6H12O	000591-78-6	86
5	2-Hexanone, 4-hydroxy-5-methyl-3...	172	C10H20O2	061141-74-0	72





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Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

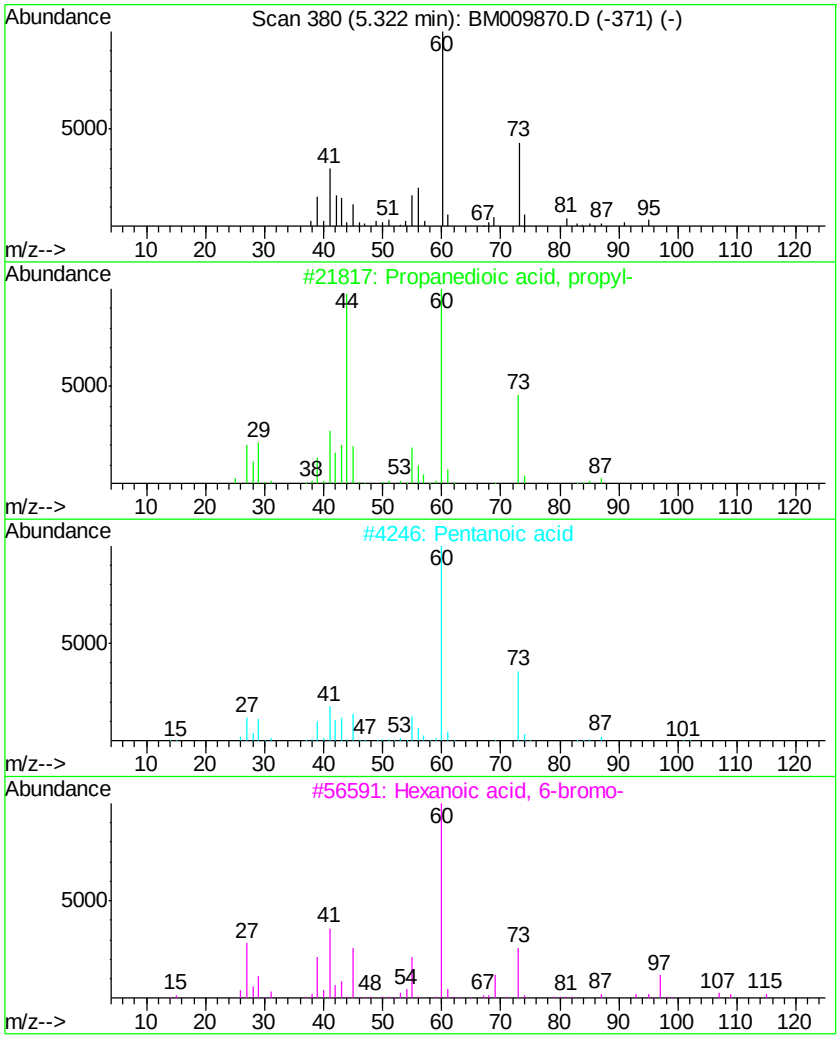
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

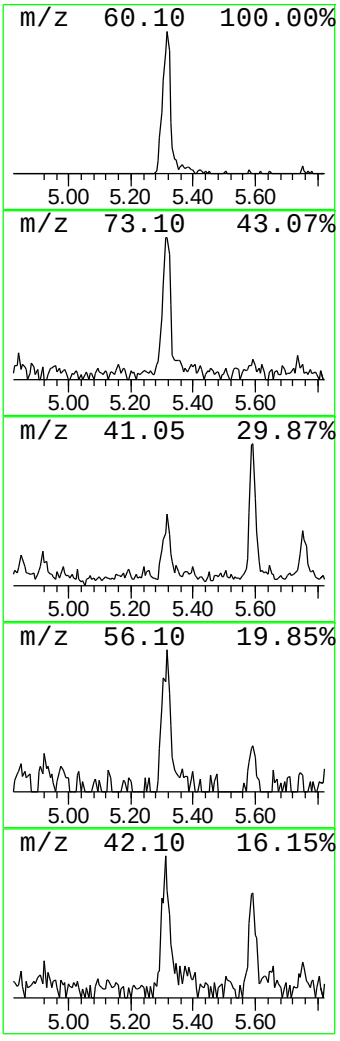
****

Peak Number 5 Propanedioic acid, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	3.74 ng/ul	146153	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	64
2	Pentanoic acid	102	C5H10O2	000109-52-4	59
3	Hexanoic acid, 6-bromo-	194	C6H11BrO2	004224-70-8	50
4	Formic acid hydrazide	60	CH4N2O	000624-84-0	43
5	Hexanoic acid	116	C6H12O2	000142-62-1	42





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

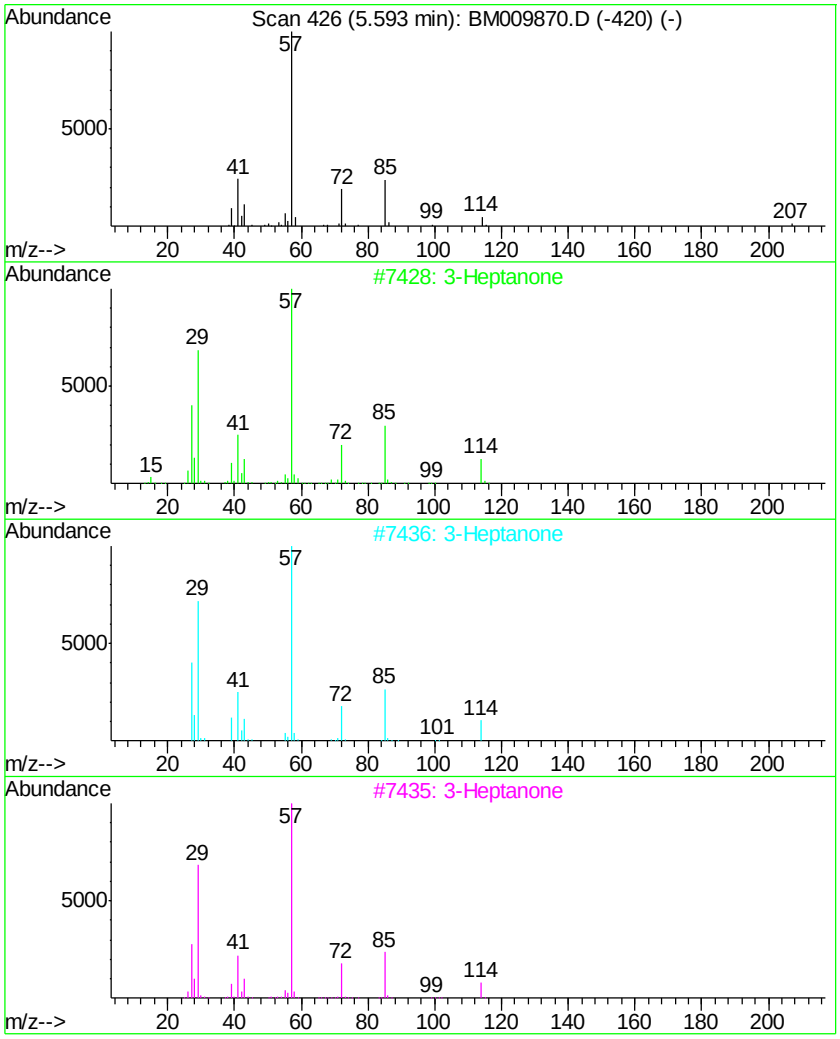
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

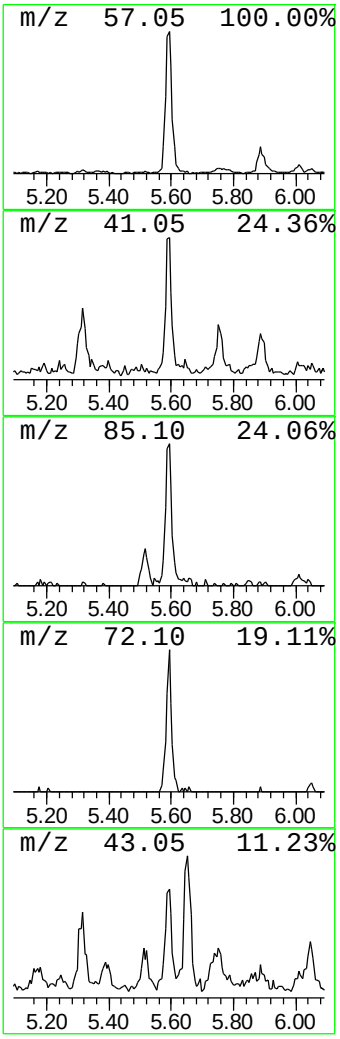
****

Peak Number 6 3-Heptanone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	5.52 ng/ul	215895	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone	114	C7H14O	000106-35-4	72
2	3-Heptanone	114	C7H14O	000106-35-4	56
3	3-Heptanone	114	C7H14O	000106-35-4	56
4	3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	53
5	3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	53





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

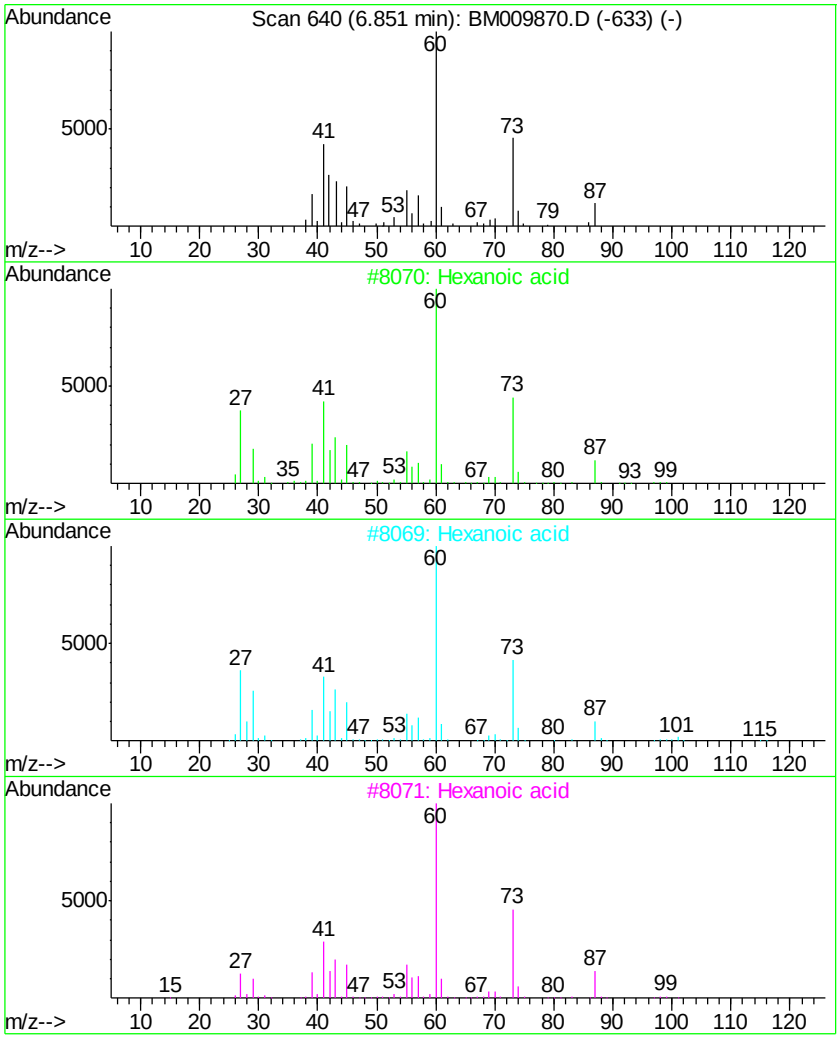
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

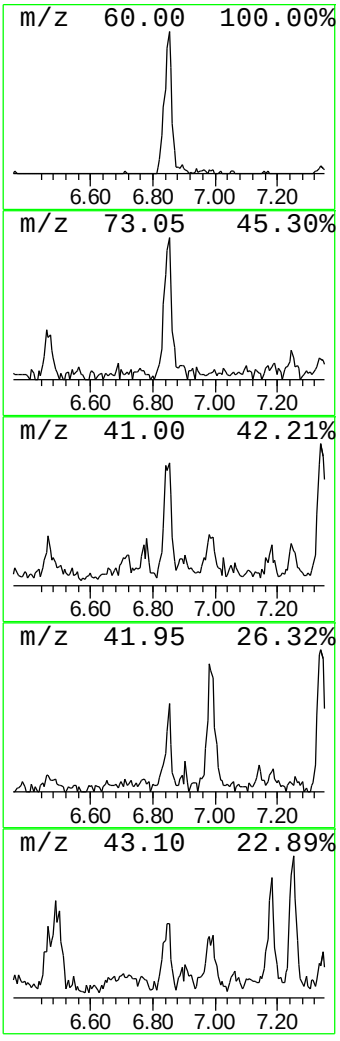
****

Peak Number 7 Hexanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.85	3.61 ng/ul	141070	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid	116	C6H12O2	000142-62-1	86
2	Hexanoic acid	116	C6H12O2	000142-62-1	64
3	Hexanoic acid	116	C6H12O2	000142-62-1	64
4	Hexanoic acid	116	C6H12O2	000142-62-1	50
5	Butanoic acid, 4-chloro-	122	C4H7ClO2	000627-00-9	50





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050317\
Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

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

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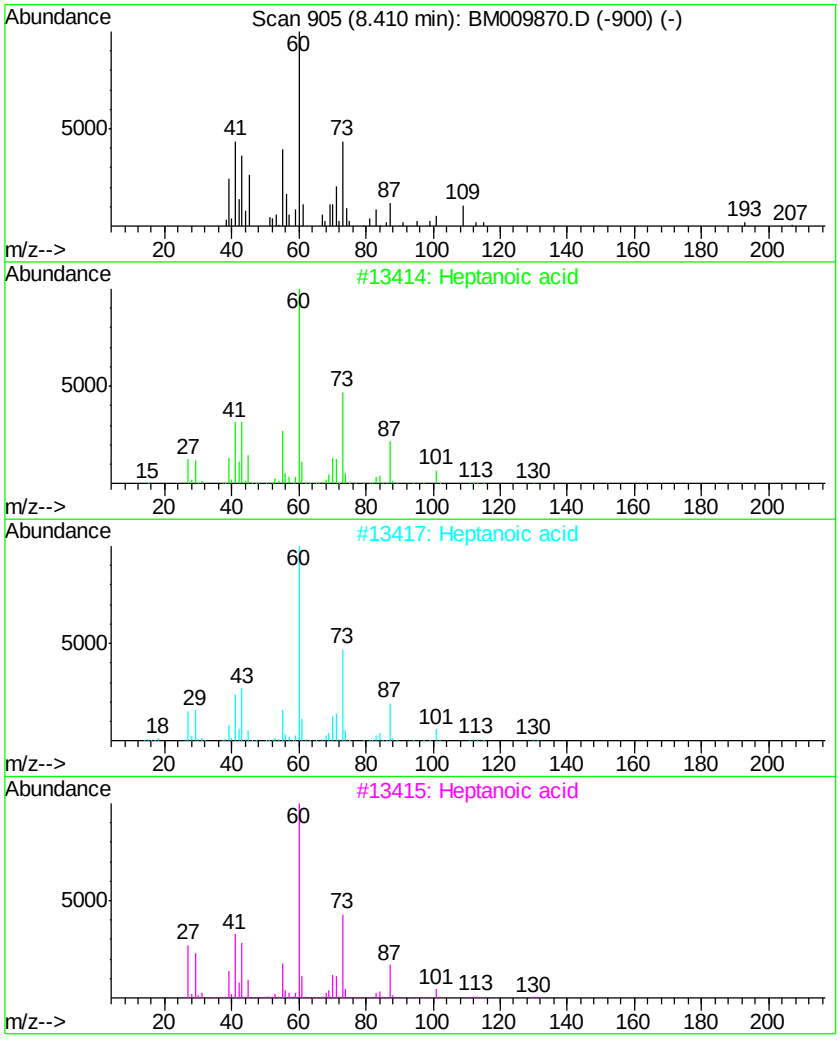
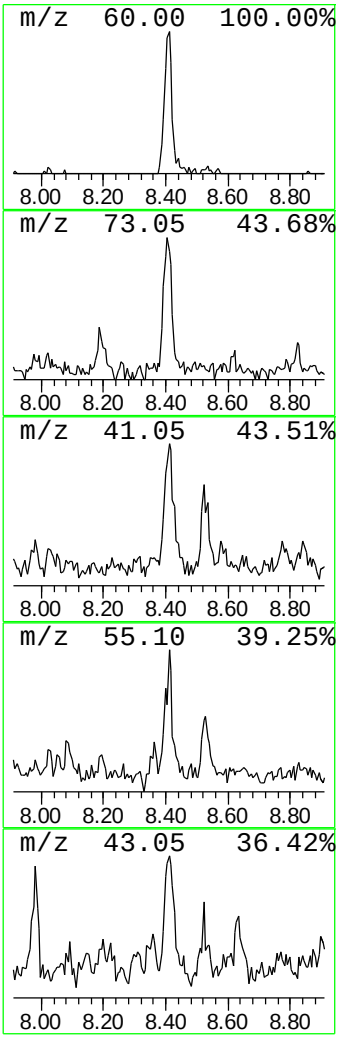
Peak Number 8 Heptanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.41	4.64 ng/ul	181611	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	64
2	Heptanoic acid	130	C7H14O2	000111-14-8	59
3	Heptanoic acid	130	C7H14O2	000111-14-8	59
4	Hexanoic acid	116	C6H12O2	000142-62-1	53
5	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	50





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Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
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TIC Library : C:\DATABASE\NIST11.L
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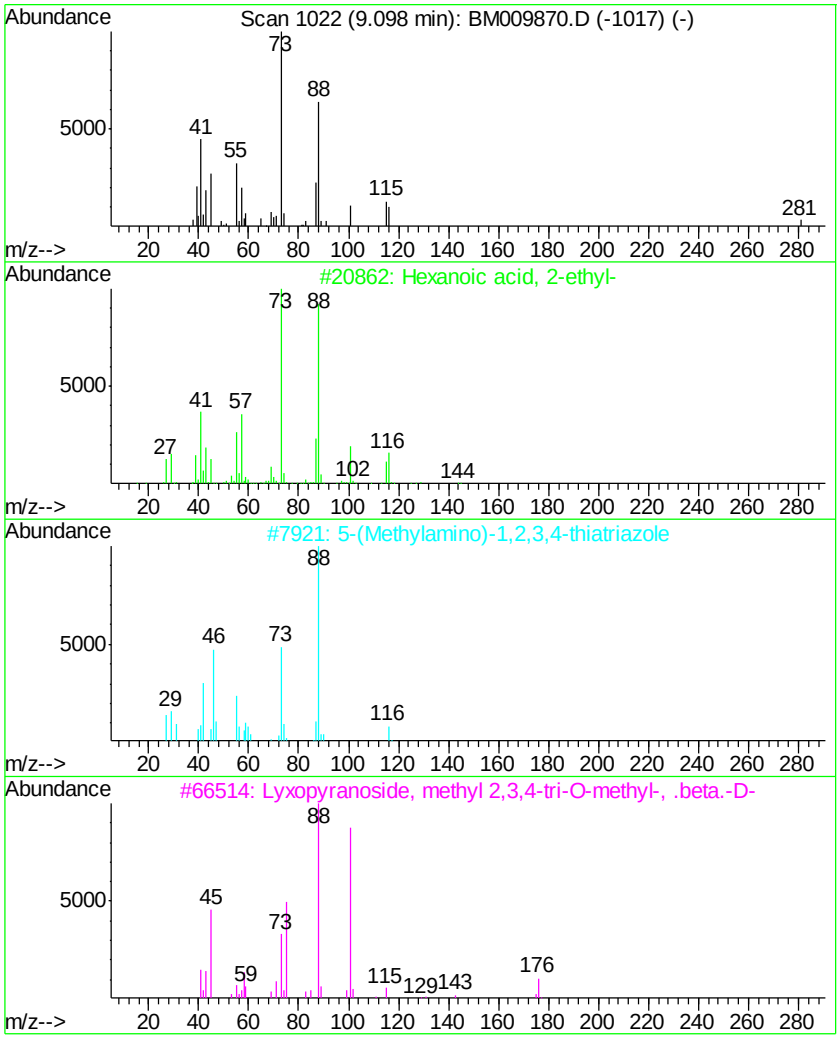
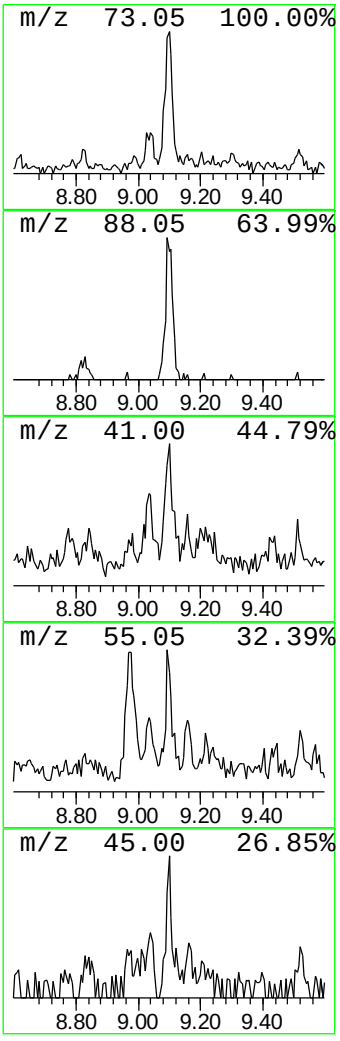
Peak Number 9 Hexanoic acid, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	2.37 ng/ul	92621	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	59
2	5-(Methylamino)-1,2,3,4-thiatria...	116	C2H4N4S	052098-72-3	43
3	Lyxopyranoside, methyl 2,3,4-tri...	206	C9H18O5	002876-84-8	33
4	Propanedioic acid, diethyl ester	160	C7H12O4	000105-53-3	33
5	2,2-Dimethylthiirane	88	C4H8S	003772-13-2	32





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Data File : BM009870.D
Acq On : 03 May 2017 12:31
Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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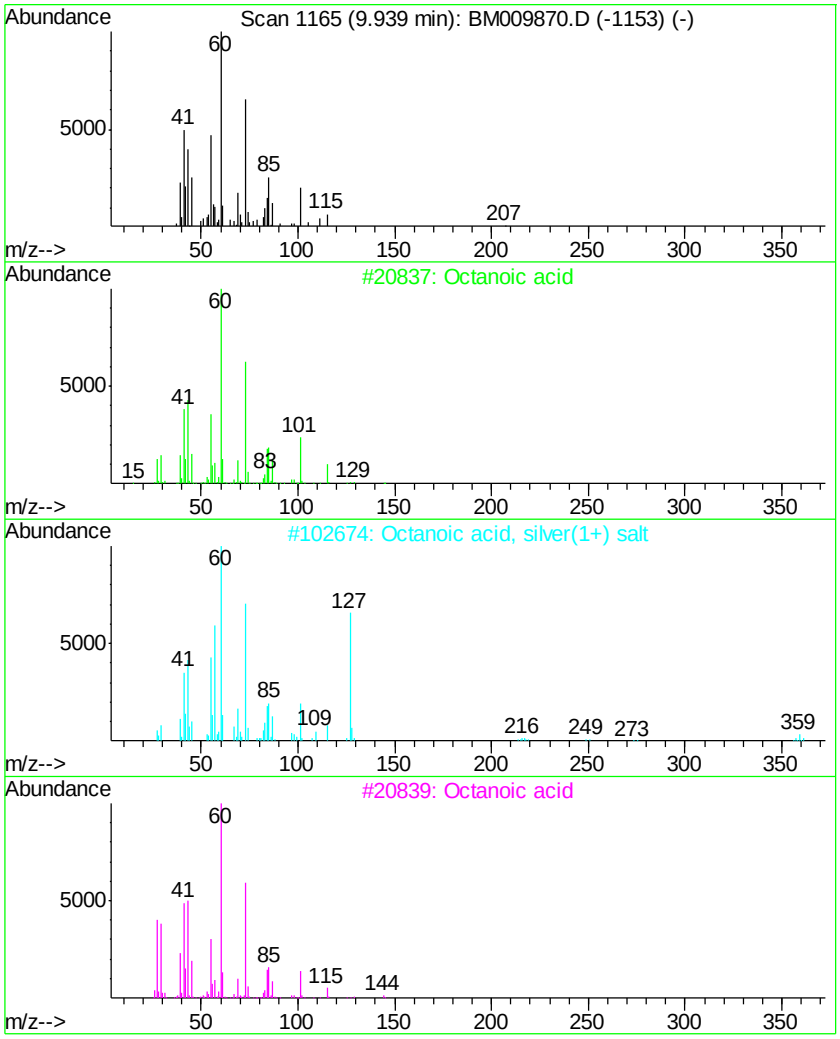
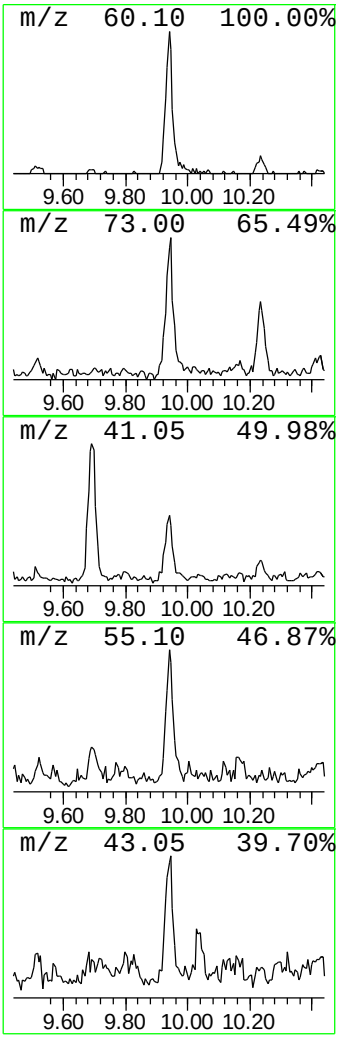
Peak Number 10 Octanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.94	3.82 ng/ul	239258	Naphthalene-d8	10.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid	144	C8H16O2	000124-07-2	64
2	Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	64
3	Octanoic acid	144	C8H16O2	000124-07-2	59
4	Hexanoic acid	116	C6H12O2	000142-62-1	59
5	Hexanoic acid	116	C6H12O2	000142-62-1	53





Library Search Compound Report

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Operator : SJ/MA
Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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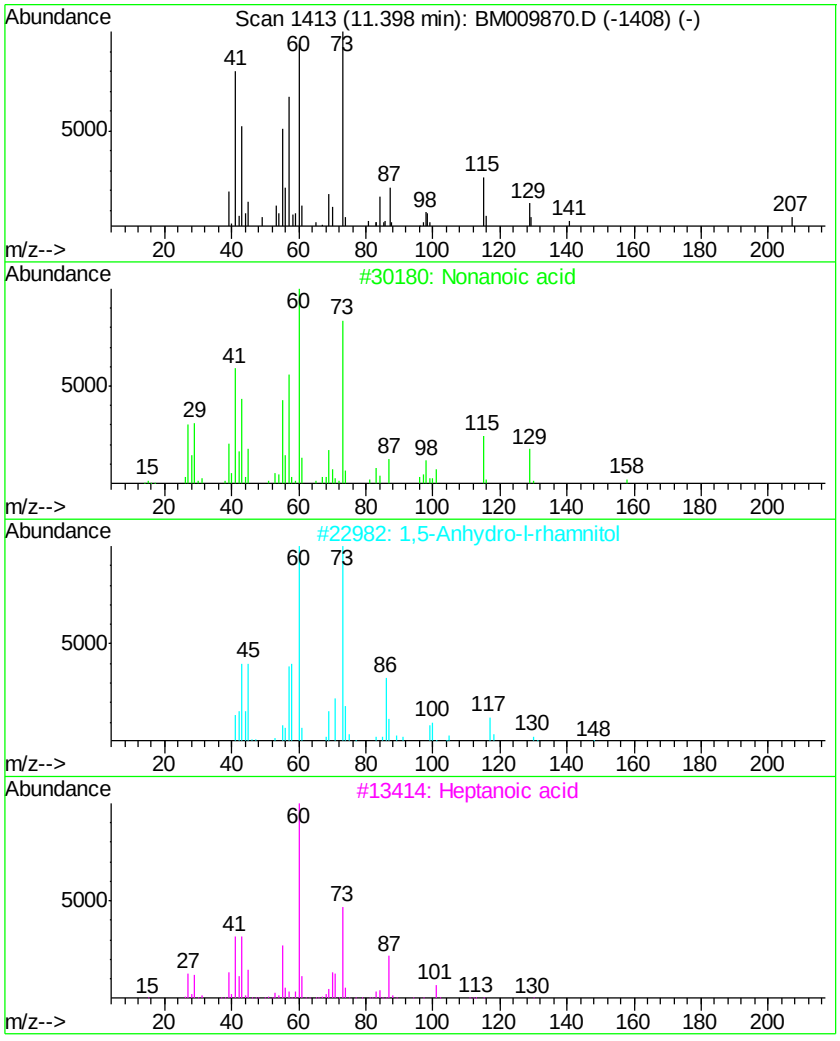
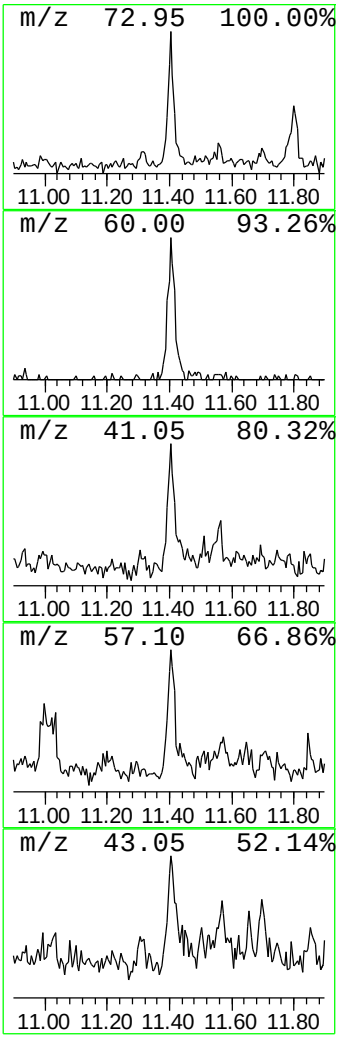
Peak Number 11 Nonanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	2.96 ng/ul	185141	Naphthalene-d8	10.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	59
2	1,5-Anhydro-l-rhamnitol	148	C6H12O4	1000129-02-4	53
3	Heptanoic acid	130	C7H14O2	000111-14-8	50
4	Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	47
5	n-Decanoic acid	172	C10H20O2	000334-48-5	42





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Sample : I2884-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : YAB41

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION

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

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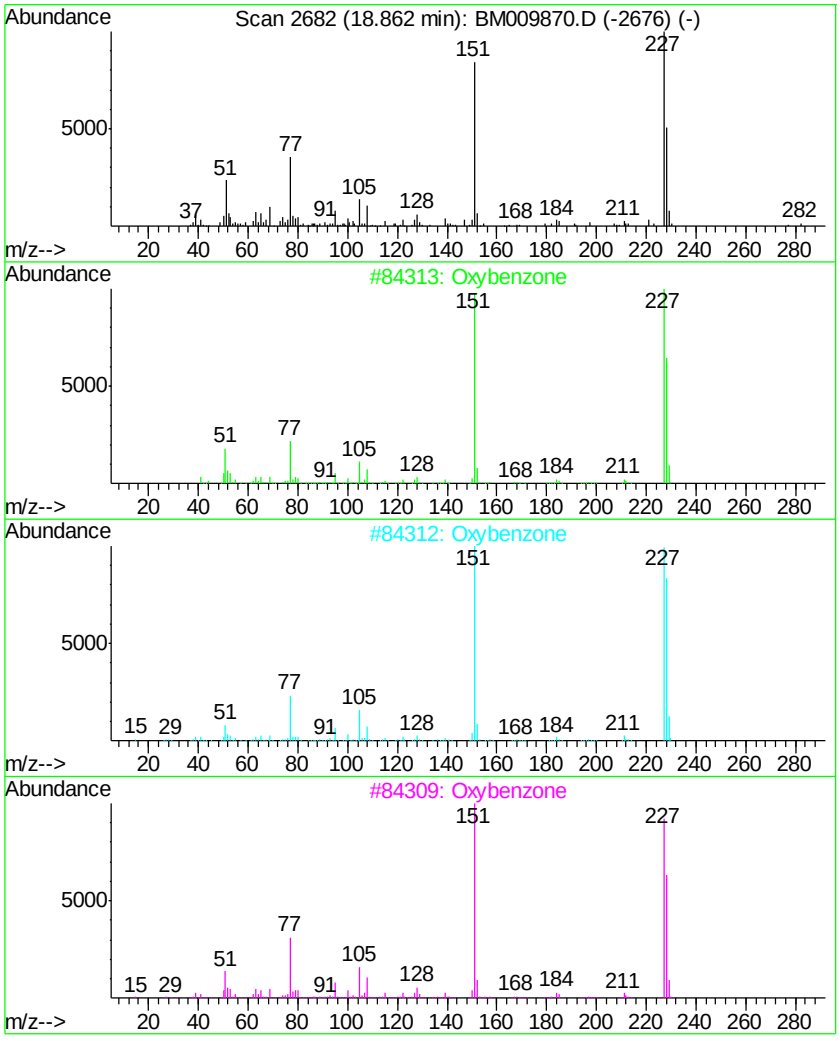
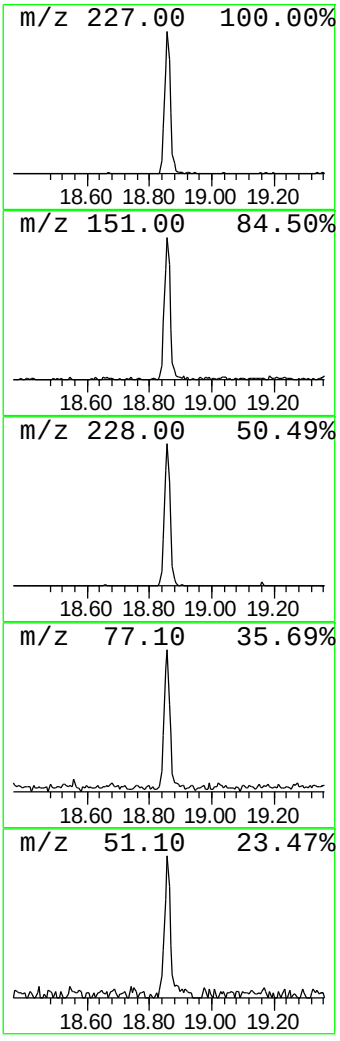
Peak Number 12 Oxybenzone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	3.60 ng/ul	331106	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxybenzone	228	C14H12O3	000131-57-7	81
2	Oxybenzone	228	C14H12O3	000131-57-7	68
3	Oxybenzone	228	C14H12O3	000131-57-7	53
4	Oxybenzone	228	C14H12O3	000131-57-7	52
5	Benzaldehyde, 3-methoxy-, oxime	151	C8H9NO2	038489-80-4	27





Tentatively Identified Compound (LSC) summary

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Operator : SJ/MA
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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
3-Pentanone	3.18	3.1	ng/ul	123251	1	7.80	782531	20.0
2-Pentanone, 3-me...	3.78	2.5	ng/ul	99275	1	7.80	782531	20.0
3-Hexanone	4.15	4.0	ng/ul	154523	1	7.80	782531	20.0
2-Hexanone	4.20	6.9	ng/ul	271036	1	7.80	782531	20.0
Propanedioic acid...	5.32	3.7	ng/ul	146153	1	7.80	782531	20.0
3-Heptanone	5.59	5.5	ng/ul	215895	1	7.80	782531	20.0
Hexanoic acid	6.85	3.6	ng/ul	141070	1	7.80	782531	20.0
Heptanoic acid	8.41	4.6	ng/ul	181611	1	7.80	782531	20.0
Hexanoic acid, 2-...	9.10	2.4	ng/ul	92621	1	7.80	782531	20.0
Octanoic acid	9.94	3.8	ng/ul	239258	2	10.60	1252450	20.0
Nonanoic acid	11.40	3.0	ng/ul	185141	2	10.60	1252450	20.0
Oxybenzone	18.86	3.6	ng/ul	331106	4	17.21	1841680	20.0