LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081318\
Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.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	5.787	471	476	494	rVB	810295	1224815	11.55%	1.455%
2	6.858	652	658	674	rVB3	237971	535659	5.05%	0.636%
3	7.022	679	686	695	rBV	790087	1248208	11.77%	1.483%
4	7.216	713	719	728	rVB	827070	1319615	12.45%	1.567%
5	7.681	791	798	805	rBV	680893	1111588	10.48%	1.320%
6	8.269	892	898	909	rBV	66508	134182	1.27%	0.159%
7	8.381	911	917	933	rVB	671469	1155029	10.89%	1.372%
8	8.769	977	983	987	rBV	80374	130037	1.23%	0.154%
9	8.828	987	993	1004	rVB	1075294	1775754	16.75%	2.109%
10	9.540	1107	1114	1124	rBV	860832	1472345	13.89%	1.749%
11	9.816	1152	1161	1172	rBV	217059	466349	4.40%	0.554%
12	10.081	1199	1206	1220	rBV	1175407	2101993	19.82%	2.497%
13	10.451	1261	1269	1277	rBV	1064595	1793028	16.91%	2.130%
14	10.816	1324	1331	1358	rBV	196864	536602	5.06%	0.637%
15	11.093	1371	1378	1391	rVB	785338	1370874	12.93%	1.628%
16	12.598	1628	1634	1644	rBV	153304	266449	2.51%	0.316%
17	13.722	1819	1825	1830	rBV	2714815	3803125	35.87%	4.517%
18	13.769	1830	1833	1840	rVB	170317	246983	2.33%	0.293%
19	13.998	1862	1872	1883	rBV	2810719	4387288	41.38%	5.211%
20	14.310	1918	1925	1936	rBV3	1759124	2733137	25.78%	3.246%
21	14.528	1956	1962	1972	rBV	229863	417969	3.94%	0.496%
22	14.757	1993	2001	2025	rBV	1019286	1777070	16.76%	2.111%
23	15.310	2087	2095	2103	rBV	3673149	5433843	51.25%	6.454%
24	15.434	2110	2116	2128	rBV	1438812	2047092	19.31%	2.431%
25	16.475	2288	2293	2303	rBV2	72829	120017	1.13%	0.143%
26	17.057	2385	2392	2399	rBV	2266017	3115065	29.38%	3.700%
27	17.157	2402	2409	2420	rVV2	4514451	6252617	58.97%	7.426%
28	17.969	2540	2547	2556	rBV2	217792	397607	3.75%	0.472%
29	18.198	2577	2586	2628	rBV	5570433	10603139	100.00%	12.594%
30	19.457	2794	2800	2808	rBV	5482022	7315758	69.00%	8.689%
31	21.051	3065	3071	3085	rBV7	57314	198547	1.87%	0.236%
32	21.257	3100	3106	3114	rBV	3009934	3830289	36.12%	4.549%
33	22.310	3281	3285	3291	rVB2	131101	167759	1.58%	0.199%
34	22.433	3302	3306	3314	rVB	493067	624508	5.89%	0.742%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081318\
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Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Title : SVOA CALIBRATION

35	22.810	3366	3370	3376	rVB	188733	255721	2.41%	0.304%
36	23.339	3452	3460	3476	rBV2	4248840	8463530	79.82%	10.052%
37	23.480	3476	3484	3494	rVB2	2128893	3934328	37.11%	4.673%
38	24.004	3568	3573	3582	rVB2	239787	440595	4.16%	0.523%
39	24.739	3691	3698	3706	rBV	209599	457559	4.32%	0.543%
40	25.592	3837	3843	3853	rVB3	131462	327745	3.09%	0.389%
41	26.598	4009	4014	4021	rVB4	78613	200186	1.89%	0.238%

Sum of corrected areas: 84194004

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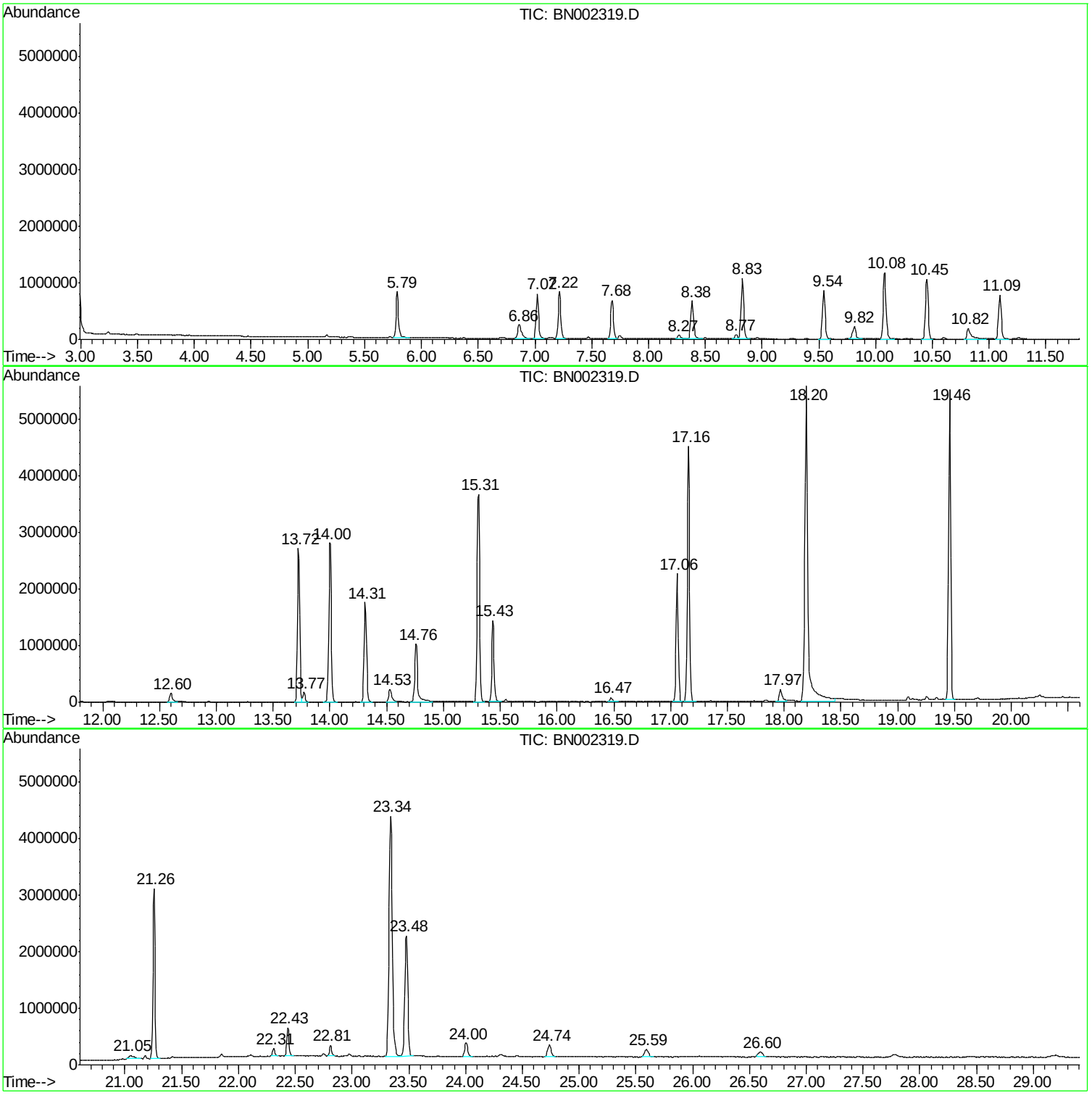
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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Data File : BN002319.D
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Operator : JU/SJ
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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

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Quant Title : SVOA CALIBRATION

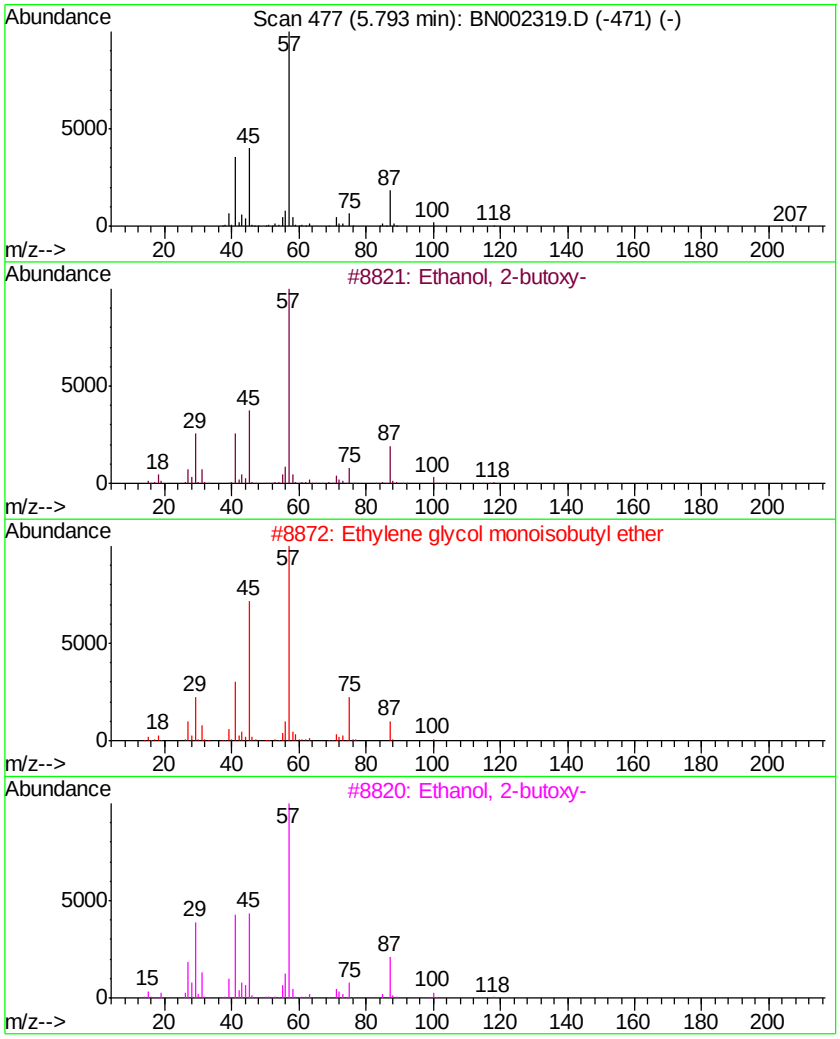
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

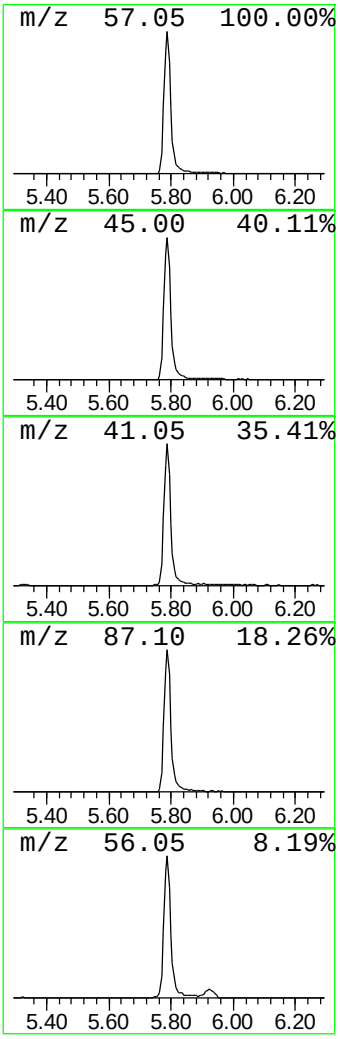
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.79	22.04 ng/ul	1224820	1,4-Dichlorobenzene-d4	7.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
2	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	53
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	50
4	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	42
5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	38





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Data File : BN002319.D
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Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

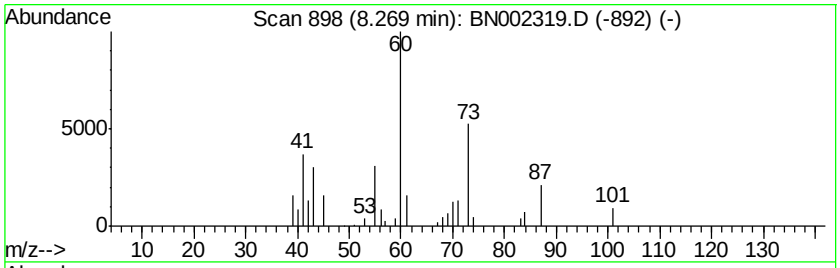
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

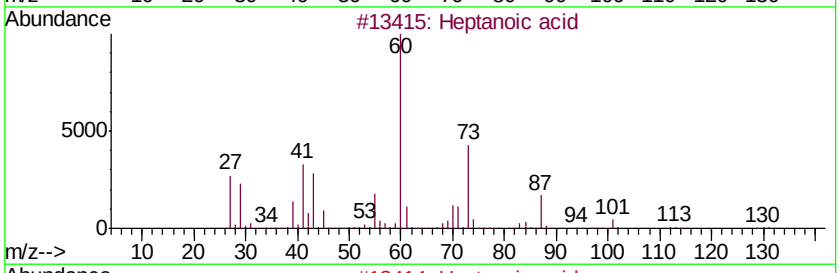
****

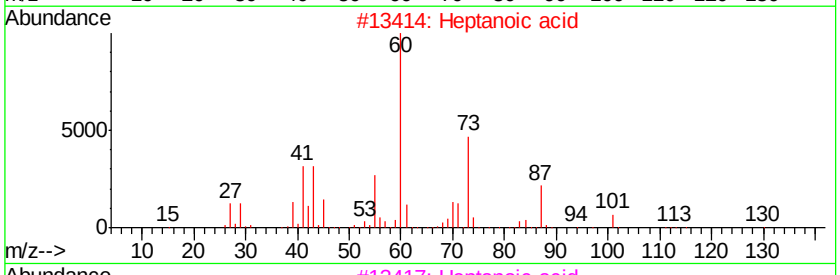
Peak Number 2 Heptanoic acid Concentration Rank 9

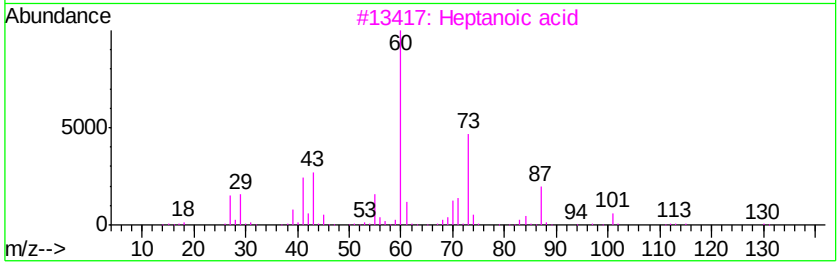
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.27	2.41 ng/ul	134182	1,4-Dichlorobenzene-d4	7.68

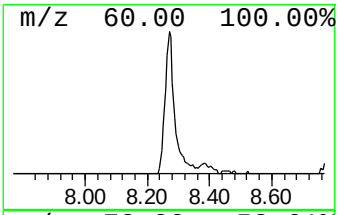
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	90
2	Heptanoic acid	130	C7H14O2	000111-14-8	90
3	Heptanoic acid	130	C7H14O2	000111-14-8	83
4	Heptanoic acid	130	C7H14O2	000111-14-8	78
5	Hexanoic acid	116	C6H12O2	000142-62-1	64

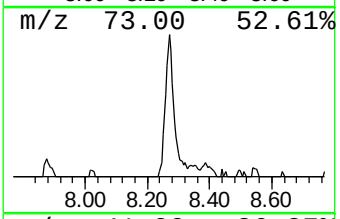


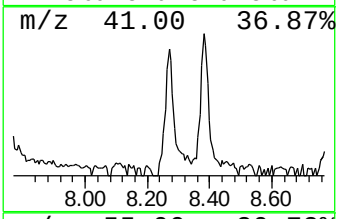


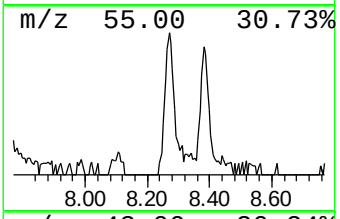


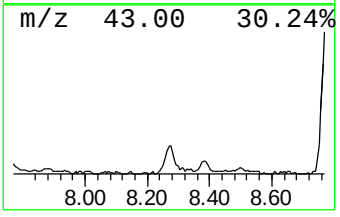












Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081318\
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Misc :
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Instrument : BNA_N
ClientSampleId : C0AC8

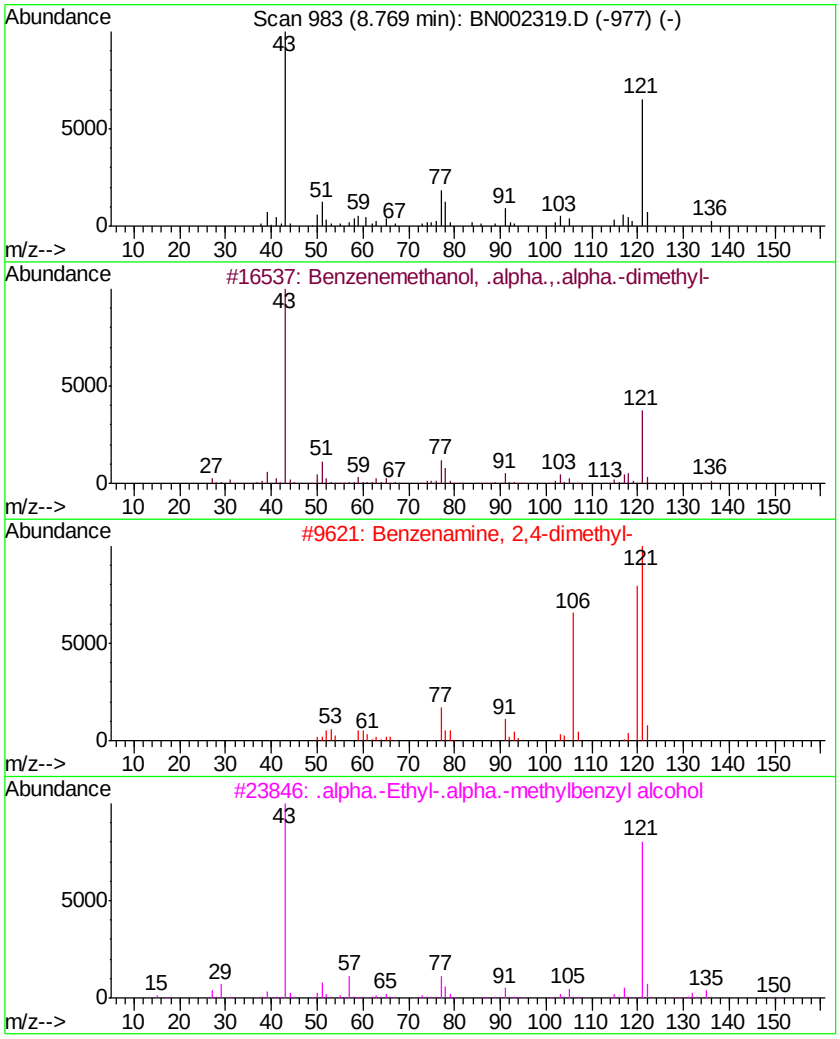
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

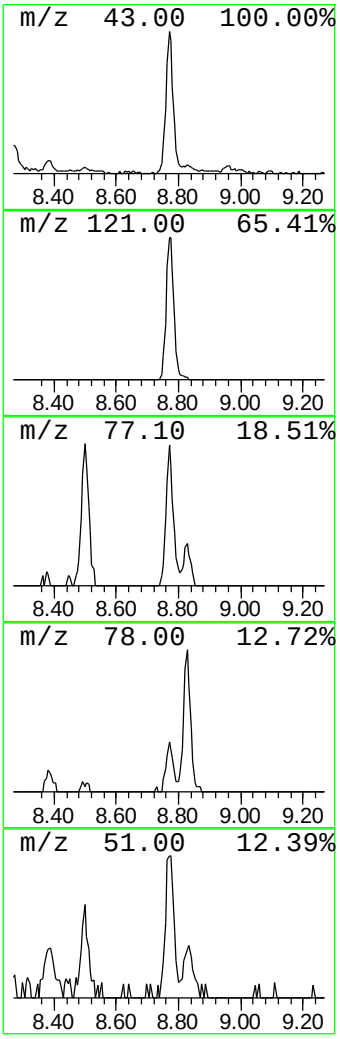
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzenemethanol, .alpha.,.a... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.77	2.34 ng/ul	130037	1,4-Dichlorobenzene-d4	7.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	68
2	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	64
3	.alpha.-Ethyl-.alpha.-methylbenz...	150	C10H14O	001565-75-9	59
4	.alpha.-Ethyl-.alpha.-methylbenz...	150	C10H14O	001565-75-9	59
5	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	58





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Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

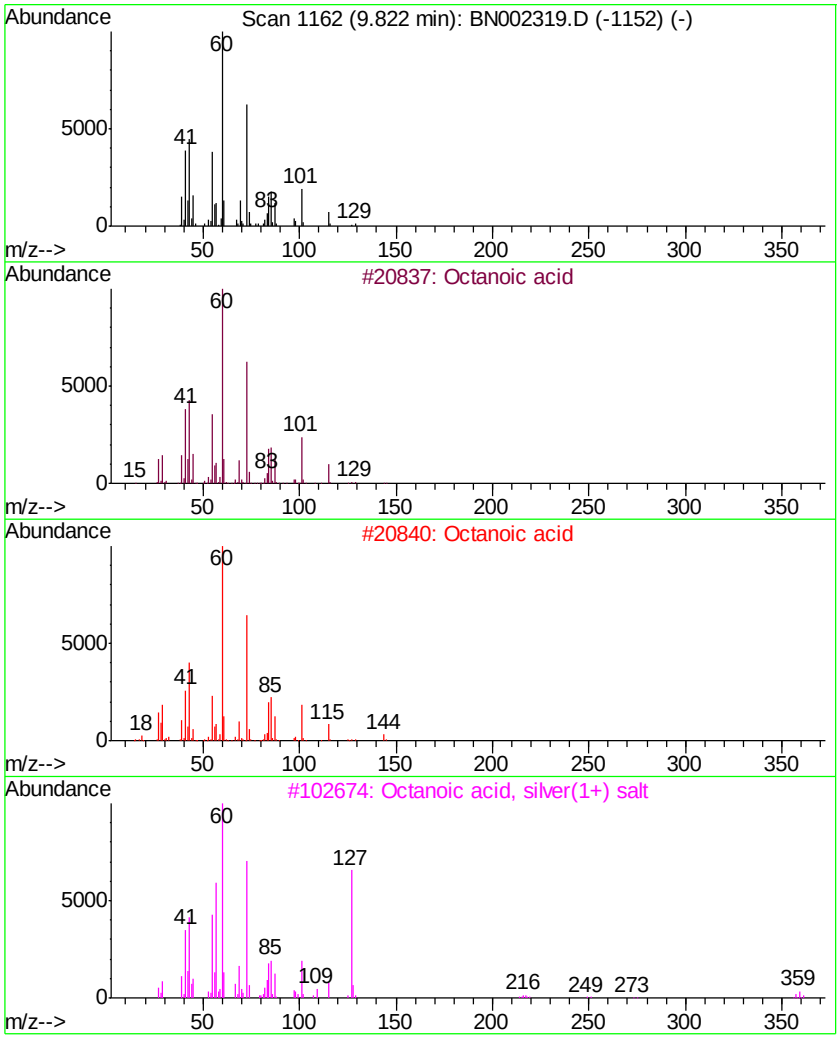
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

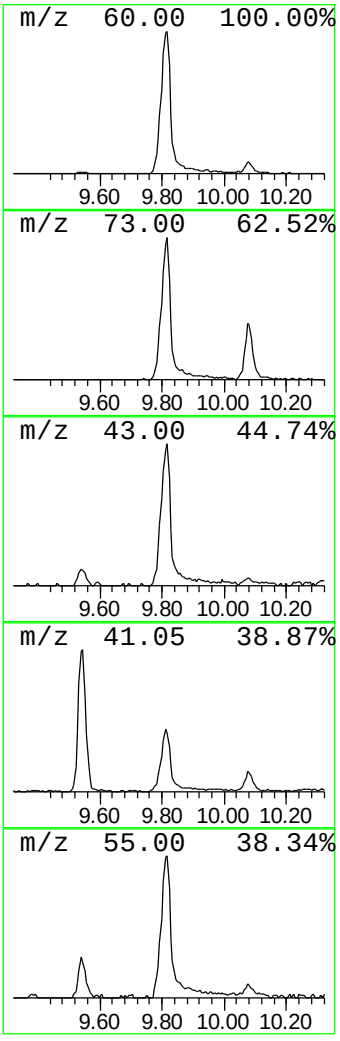
****

Peak Number 4 Octanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	5.20 ng/ul	466349	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid	144	C8H16O2	000124-07-2	91
2	Octanoic acid	144	C8H16O2	000124-07-2	86
3	Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	83
4	Octanoic acid	144	C8H16O2	000124-07-2	78
5	Hexanoic acid	116	C6H12O2	000142-62-1	59





Library Search Compound Report

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Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

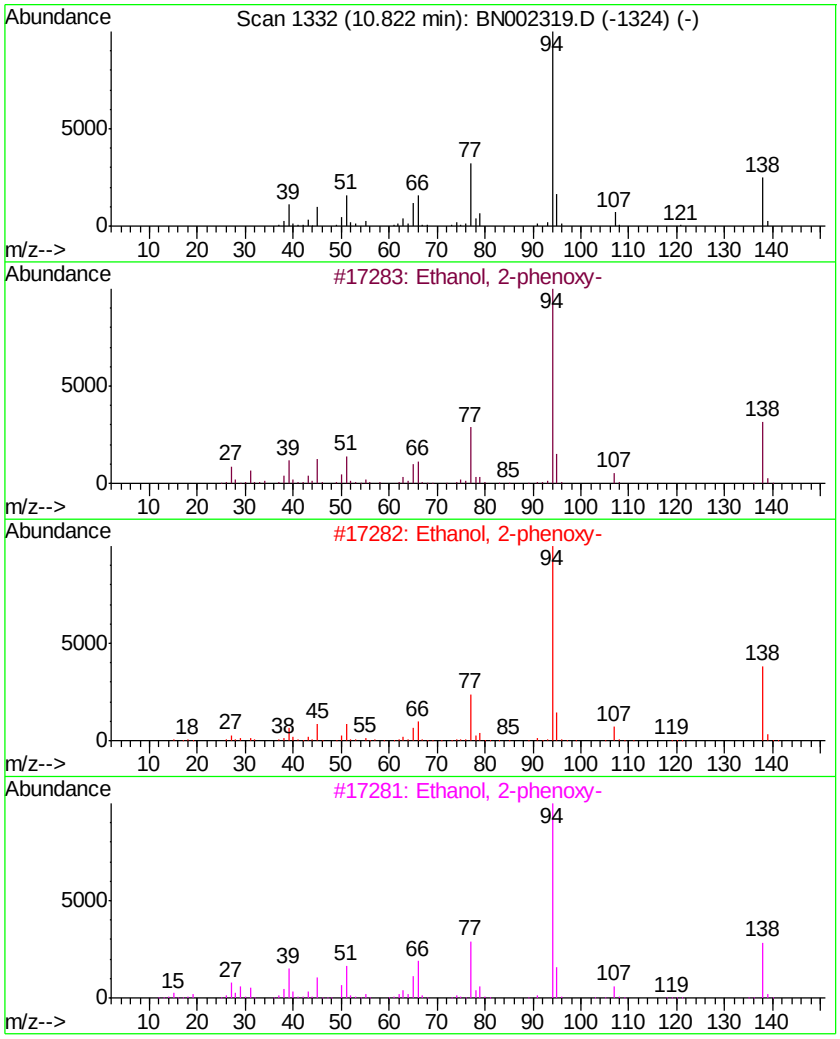
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

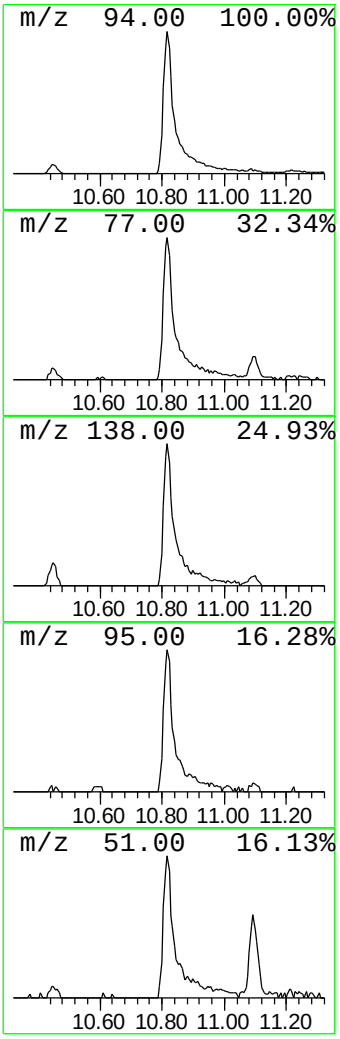
****

Peak Number 5 Ethanol, 2-phenoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.82	5.99 ng/ul	536602	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	95
2	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
4	Methane, isopropoxyphenoxy-	166	C10H14O2	000828-12-6	53
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	53





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Instrument : BNA_N
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Quant Title : SVOA CALIBRATION

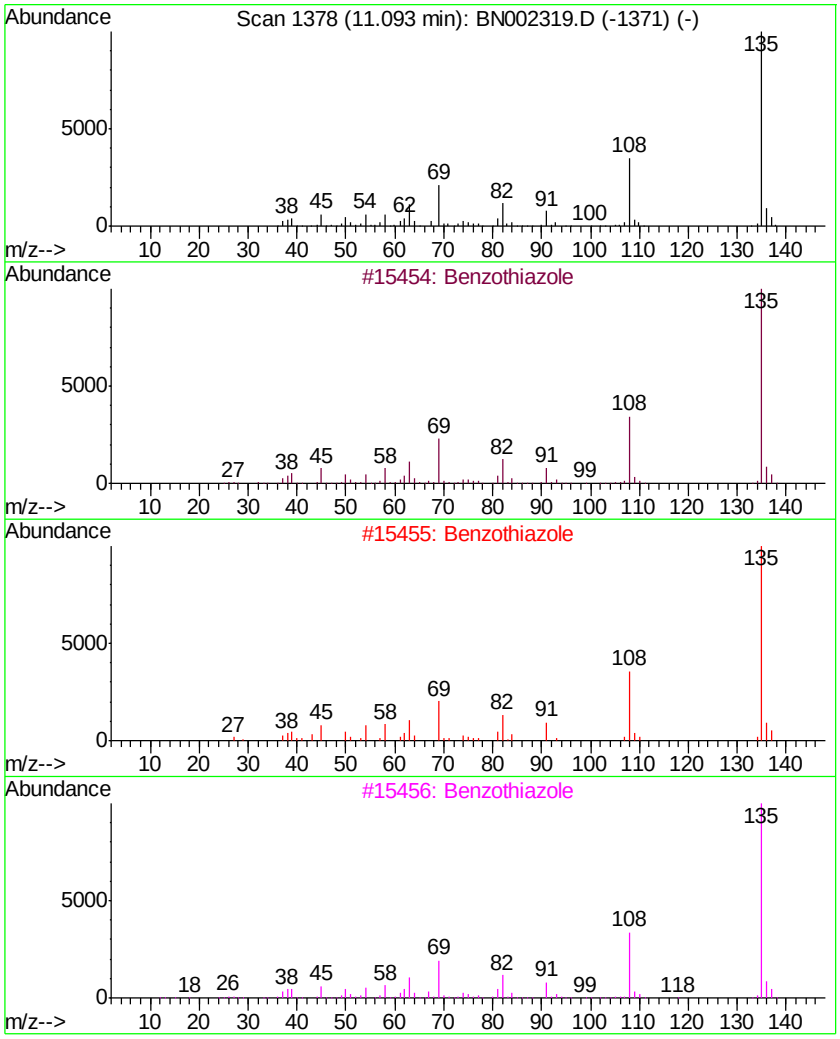
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

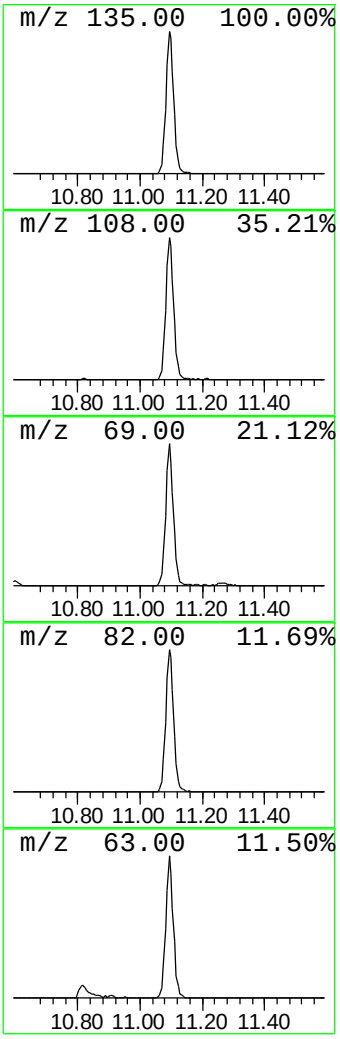
****

Peak Number 6 Benzothiazole Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	15.29 ng/ul	1370870	Naphthalene-d8	10.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzothiazole	135	C7H5NS	000095-16-9	97
2	Benzothiazole	135	C7H5NS	000095-16-9	97
3	Benzothiazole	135	C7H5NS	000095-16-9	95
4	1,2-Benzisothiazole	135	C7H5NS	000272-16-2	90
5	1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	59





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ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

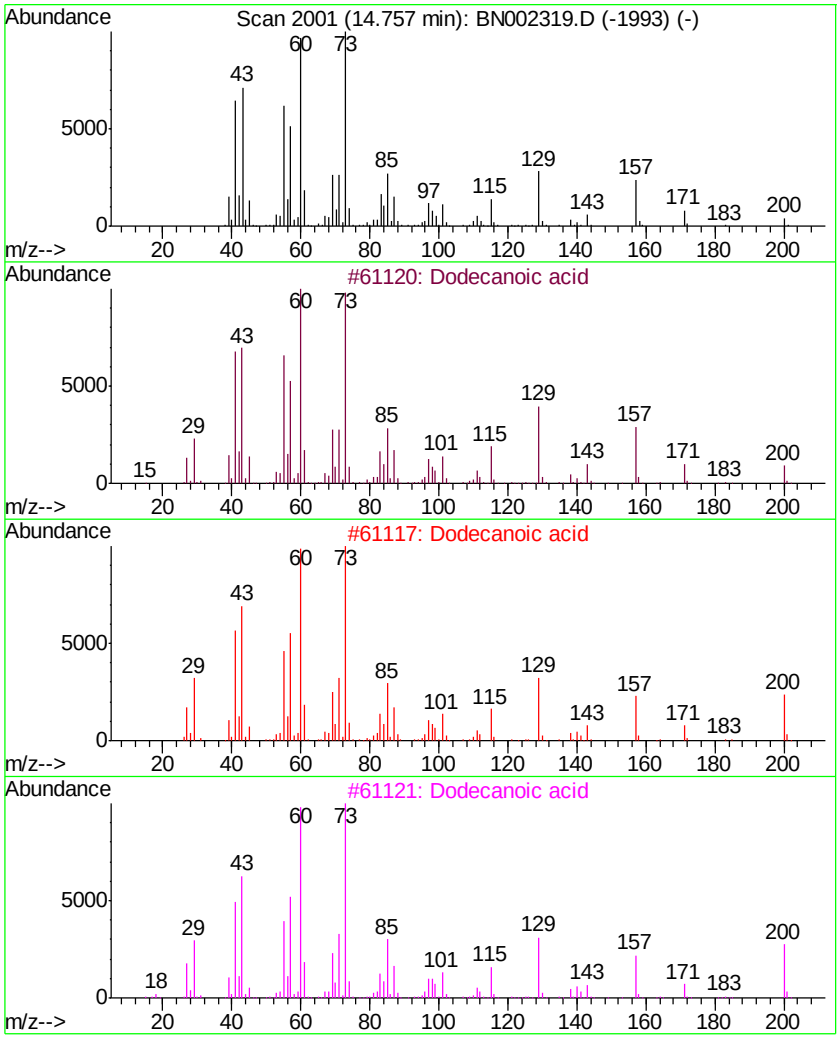
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

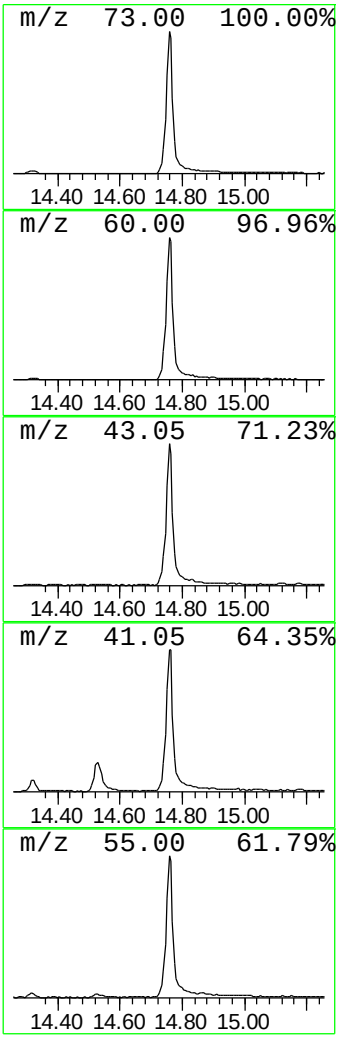
****

Peak Number 7 Dodecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	13.00 ng/ul	1777070	Acenaphthene-d10	14.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid	200	C12H24O2	000143-07-7	98
2	Dodecanoic acid	200	C12H24O2	000143-07-7	91
3	Dodecanoic acid	200	C12H24O2	000143-07-7	91
4	Undecanoic acid	186	C11H22O2	000112-37-8	80
5	n-Decanoic acid	172	C10H20O2	000334-48-5	70





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Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

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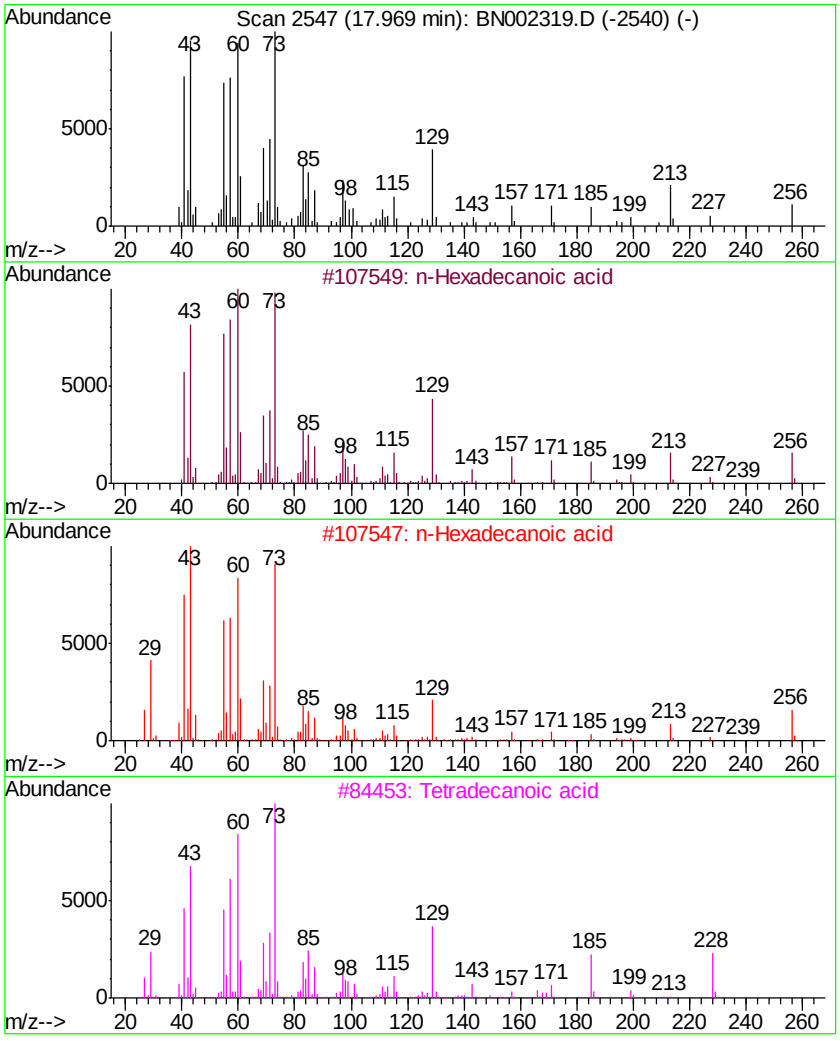
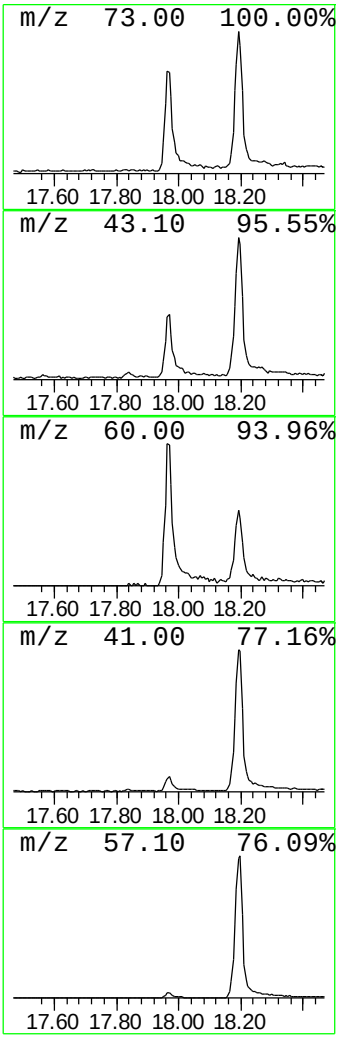
Peak Number 8 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	2.55 ng/ul	397607	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4	Tridecanoic acid	214	C13H26O2	000638-53-9	68
5	Tridecanoic acid	214	C13H26O2	000638-53-9	68





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Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

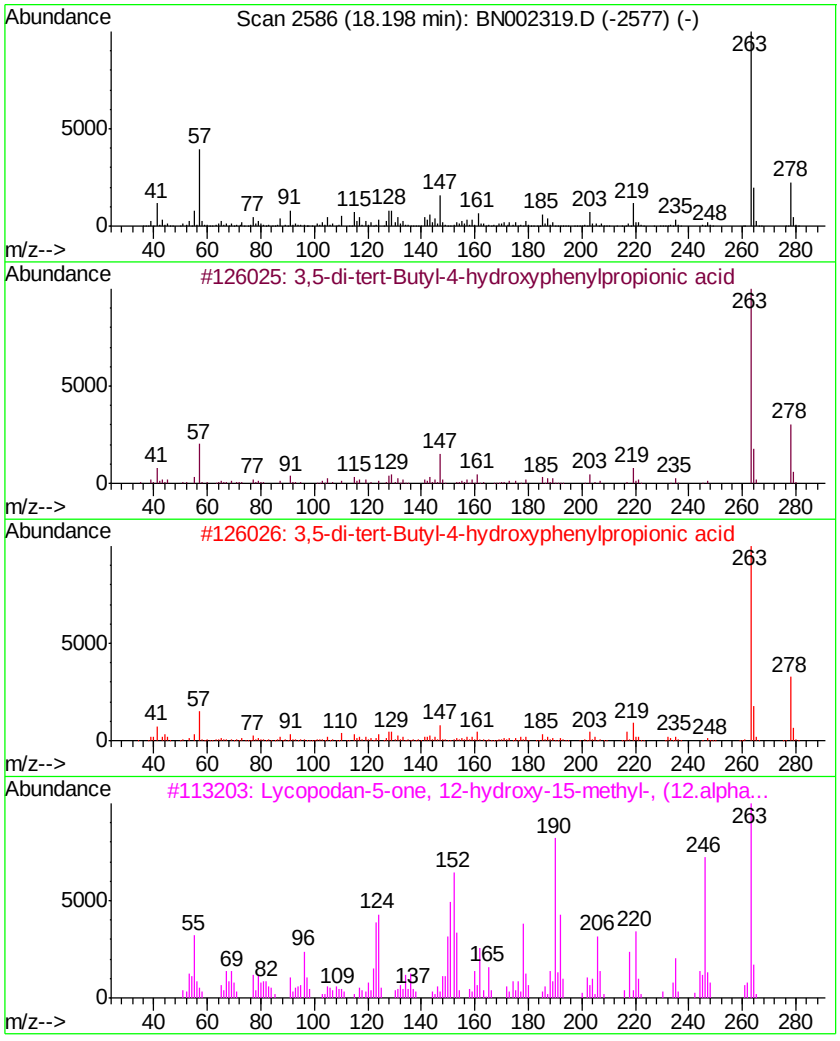
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

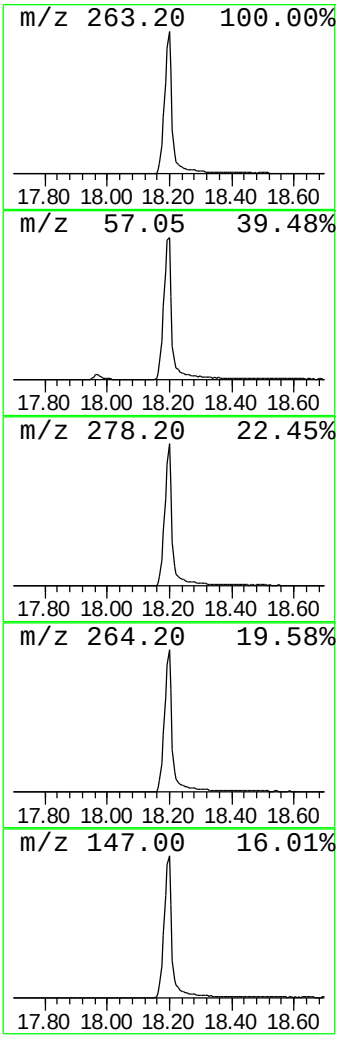
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.20	68.08 ng/ul	10603100	Phenanthrene-d10	17.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	95
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3	Lycopodan-5-one, 12-hydroxy-15-m...	263	C16H25NO2	021061-90-5	91
4	2,2-Bis[4'-cyanooxyphenyl]propane	278	C17H14N2O2	1000322-13-3	72
5	2-(4-Methoxy-phenyl)-1H-1,3a,8-t...	263	C16H13N3O	036289-23-3	53





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Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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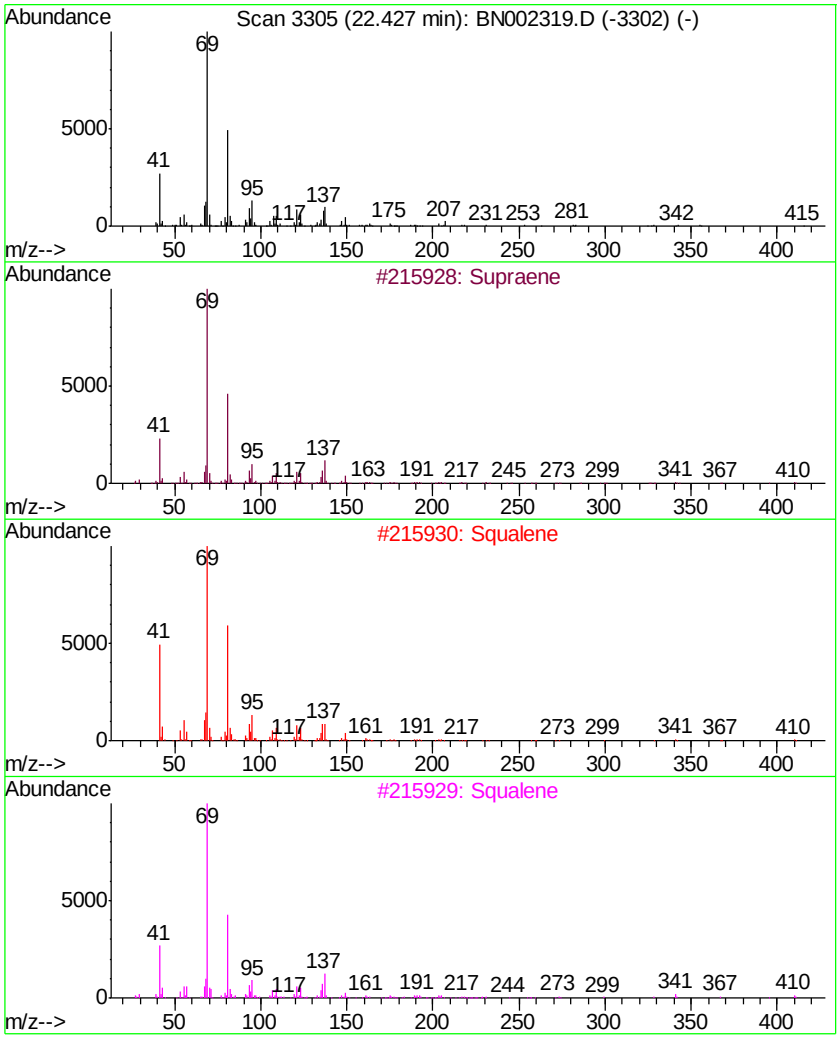
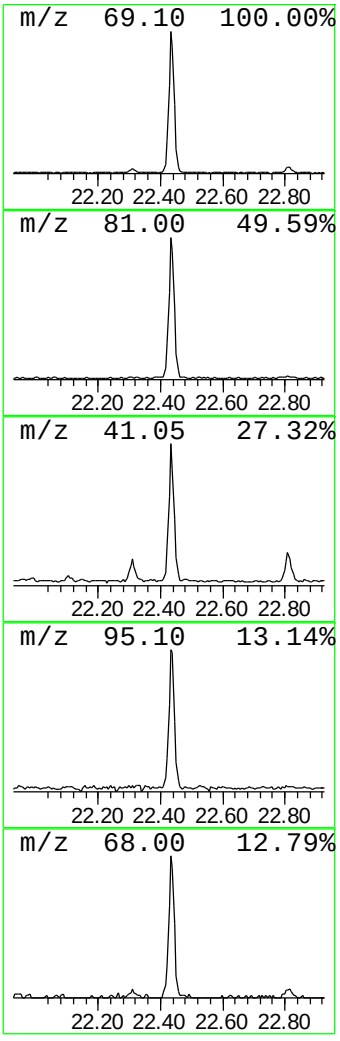
Peak Number 10 Supraene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.43	3.17 ng/ul	624508	Perylene-d12	23.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	91
2	Squalene	410	C30H50	000111-02-4	91
3	Squalene	410	C30H50	000111-02-4	90
4	Farnesol isomer a	222	C15H26O	1000108-92-4	64
5	2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	52





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Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

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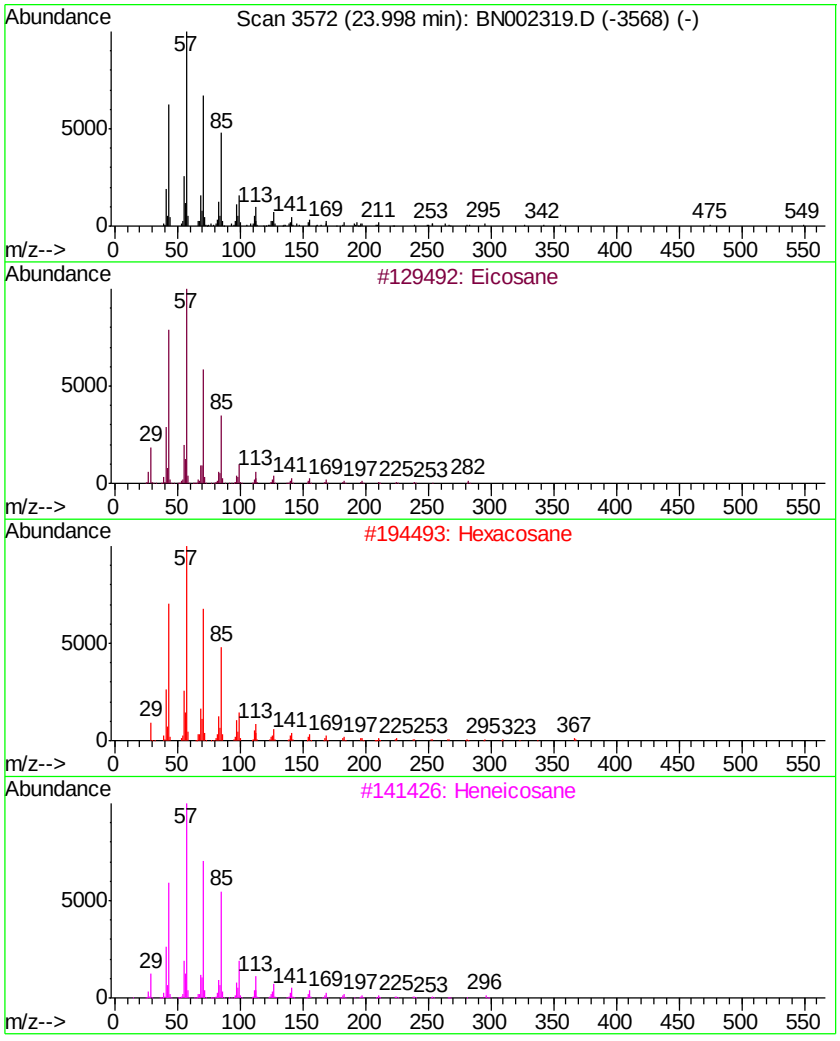
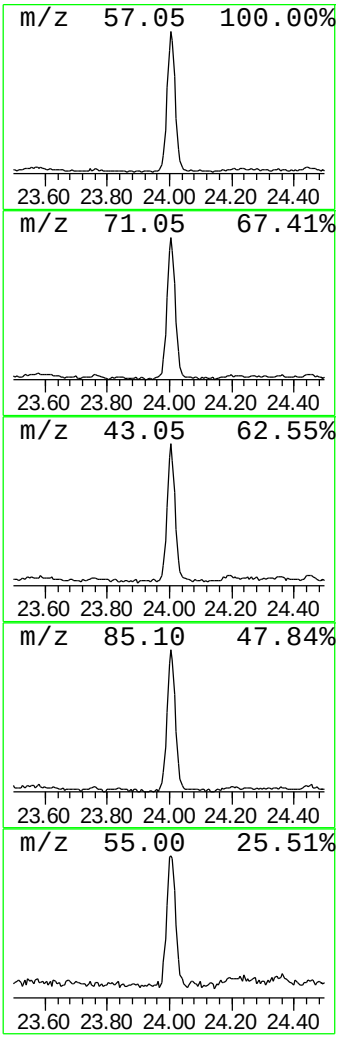
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	2.24 ng/ul	440595	Perylene-d12	23.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Hexacosane	366	C26H54	000630-01-3	93
3	Heneicosane	296	C21H44	000629-94-7	90
4	Heptacosane	380	C27H56	000593-49-7	90
5	Octacosane	394	C28H58	000630-02-4	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081318\
Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

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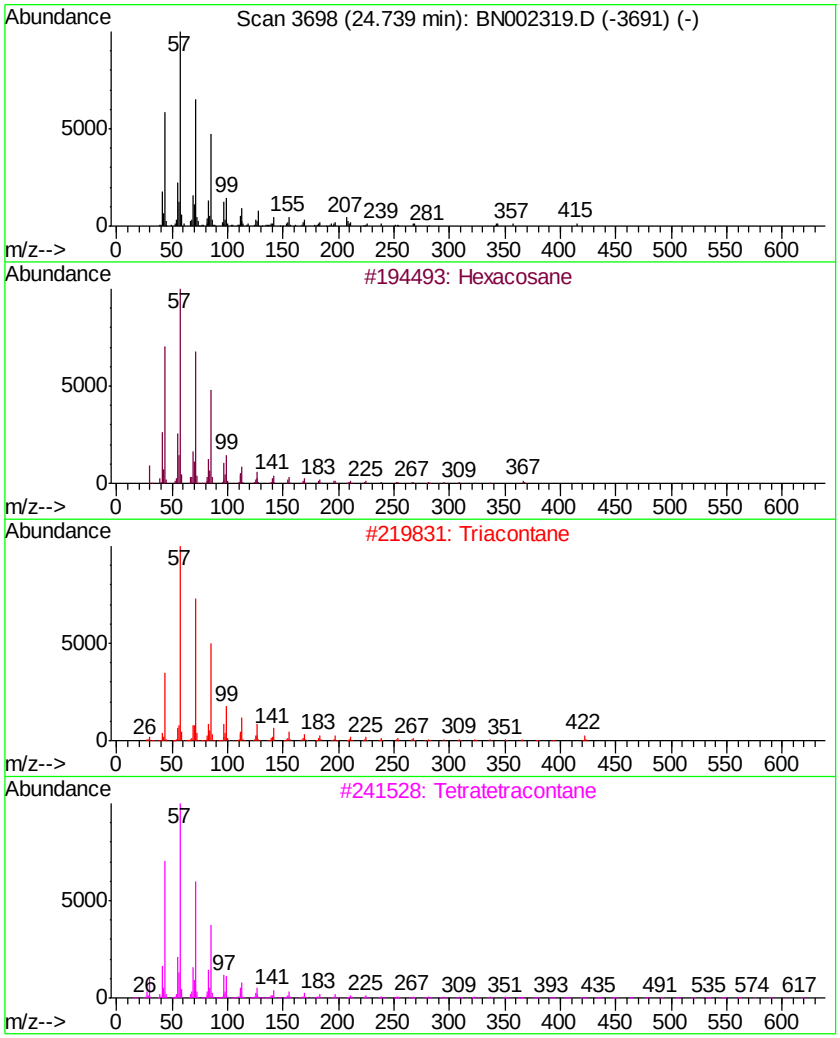
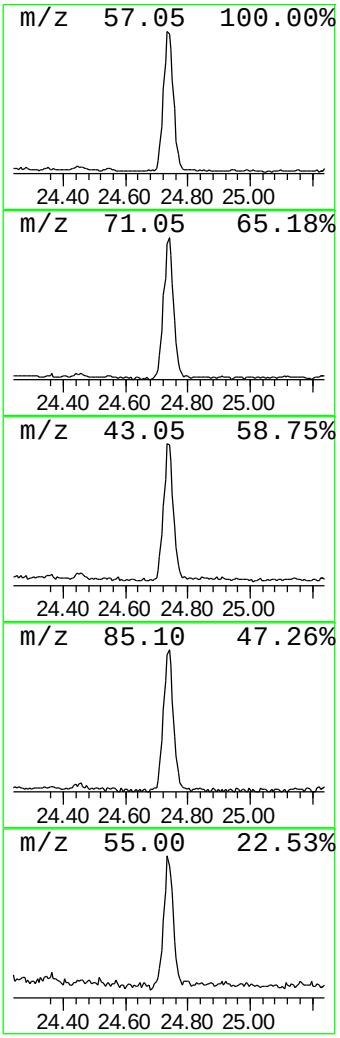
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.74	2.33 ng/ul	457559	Perylene-d12	23.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	93
2	Triacontane	422	C30H62	000638-68-6	90
3	Tetratetracontane	619	C44H90	007098-22-8	90
4	Heneicosane	296	C21H44	000629-94-7	90
5	Octacosane	394	C28H58	000630-02-4	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081318\
Data File : BN002319.D
Acq On : 13 Aug 2018 22:53
Operator : JU/SJ
Sample : J4422-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0AC8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.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
Ethanol, 2-butoxy-	5.79	22.0	ng/ul	1224820	1	7.68	1111590	20.0
Heptanoic acid	8.27	2.4	ng/ul	134182	1	7.68	1111590	20.0
Benzenemethanol, ...	8.77	2.3	ng/ul	130037	1	7.68	1111590	20.0
Octanoic acid	9.82	5.2	ng/ul	466349	2	10.45	1793030	20.0
Ethanol, 2-phenoxy-	10.82	6.0	ng/ul	536602	2	10.45	1793030	20.0
Benzothiazole	11.09	15.3	ng/ul	1370870	2	10.45	1793030	20.0
Dodecanoic acid	14.76	13.0	ng/ul	1777070	3	14.31	2733140	20.0
n-Hexadecanoic acid	17.97	2.5	ng/ul	397607	4	17.06	3115070	20.0
3,5-di-tert-Butyl...	18.20	68.1	ng/ul	10603100	4	17.06	3115070	20.0
Supraene	22.43	3.2	ng/ul	624508	6	23.48	3934330	20.0
(DEL) Alkane: Str...	24.00	2.2	ng/ul	440595	6	23.48	3934330	20.0
(DEL) Alkane: Str...	24.74	2.3	ng/ul	457559	6	23.48	3934330	20.0