LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_G\METHODS\8270-BG061719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.115	3	4	17	rVB	527153	625075	2.73%	1.082%
2	5.676	432	440	452	rVB	192780	407942	1.78%	0.706%
3	6.187	505	527	531	rBV	5719089	22895164	100.00%	39.638%
4	7.216	694	702	715	rBV	158728	311996	1.36%	0.540%
5	7.586	756	765	777	rBV	515021	904260	3.95%	1.566%
6	8.056	837	845	850	rBV	161965	279676	1.22%	0.484%
7	8.379	891	900	910	rBV	537521	920434	4.02%	1.594%
8	9.108	1001	1024	1038	rBV2	336461	1451131	6.34%	2.512%
9	9.237	1039	1046	1054	rVV	424375	763551	3.33%	1.322%
10	10.876	1317	1325	1339	rBV	187200	344694	1.51%	0.597%
11	10.999	1339	1346	1372	rBV2	181725	727994	3.18%	1.260%
12	13.314	1729	1740	1748	rVB	1253606	1895622	8.28%	3.282%
13	13.632	1788	1794	1806	rBV2	295133	586885	2.56%	1.016%
14	14.695	1970	1975	1980	rVB2	300831	508322	2.22%	0.880%
15	15.905	2175	2181	2185	rBV	605827	1049355	4.58%	1.817%
16	16.193	2225	2230	2237	rBV2	1331433	2763969	12.07%	4.785%
17	16.393	2258	2264	2268	rBV	2387369	3850085	16.82%	6.666%
18	17.222	2400	2405	2409	rBV	2689860	3951997	17.26%	6.842%
19	17.827	2505	2508	2512	rVB	1171132	1458562	6.37%	2.525%
20	18.273	2581	2584	2590	rVB	1358118	1909034	8.34%	3.305%
21	19.924	2862	2865	2869	rBV	1535566	1868919	8.16%	3.236%
22	20.059	2885	2888	2892	rVB	2409991	2902459	12.68%	5.025%
23	21.305	3096	3100	3112	rVB	1373473	2440815	10.66%	4.226%
24	22.938	3372	3378	3390	rVB	649563	1395098	6.09%	2.415%
25	25.018	3726	3732	3745	rVB2	217219	624816	2.73%	1.082%
26	25.341	3780	3787	3807	rVB2	284256	923391	4.03%	1.599%

Sum of corrected areas: 57761246

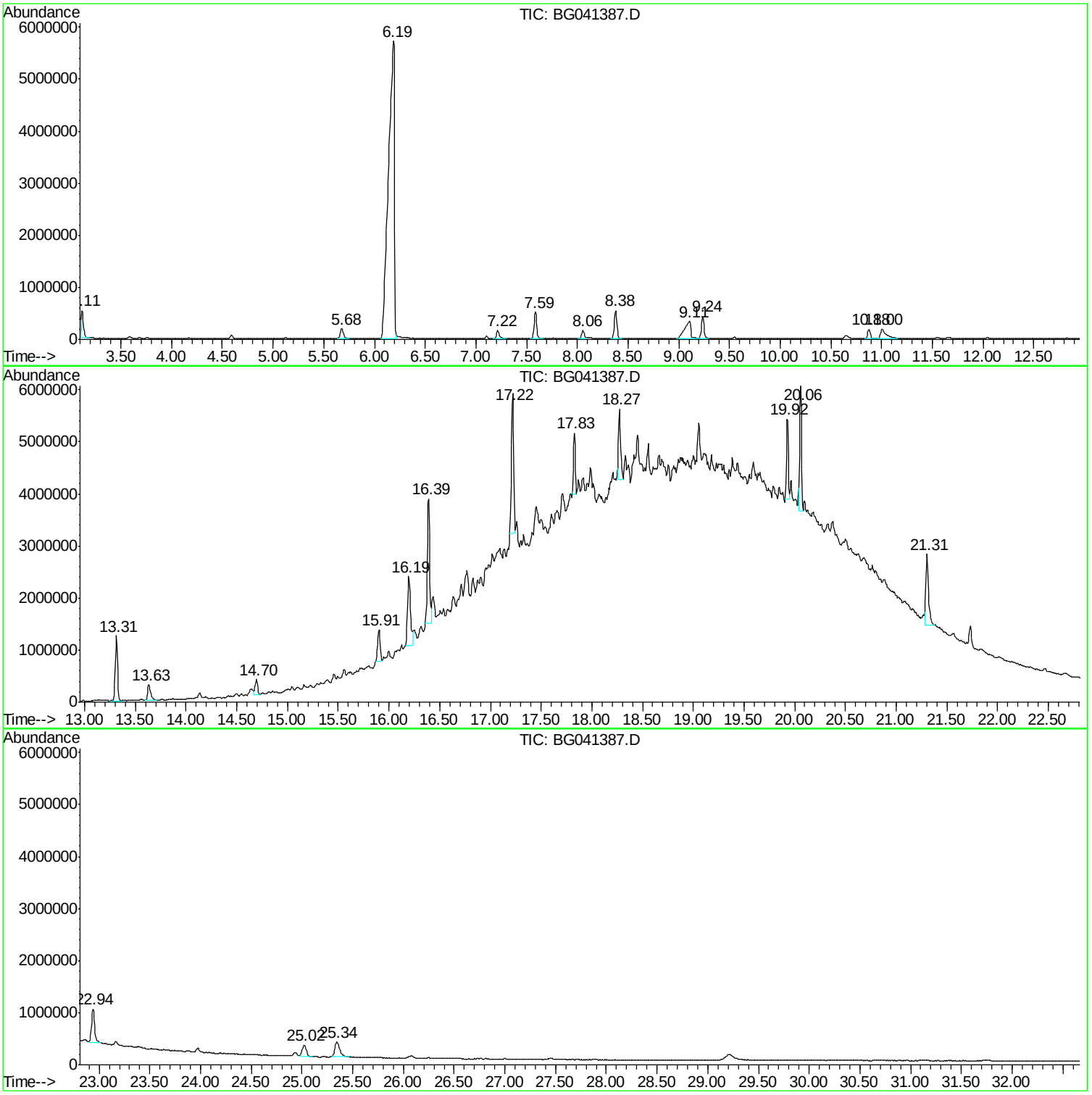
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

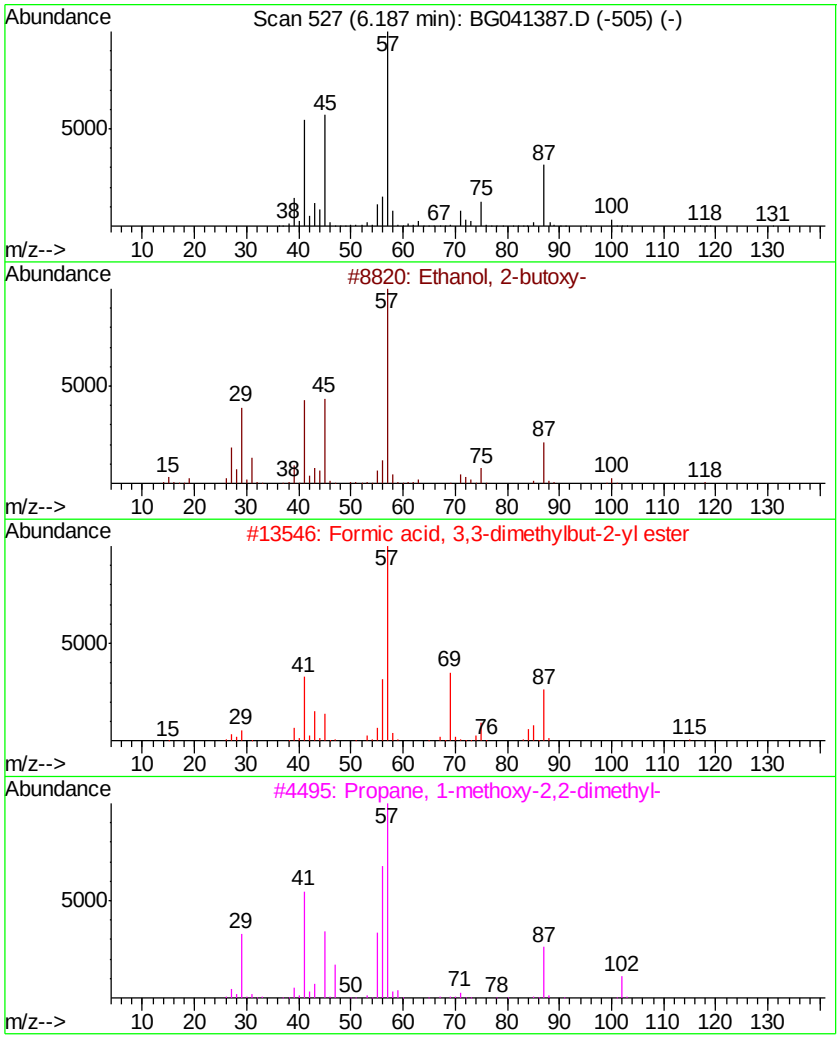
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

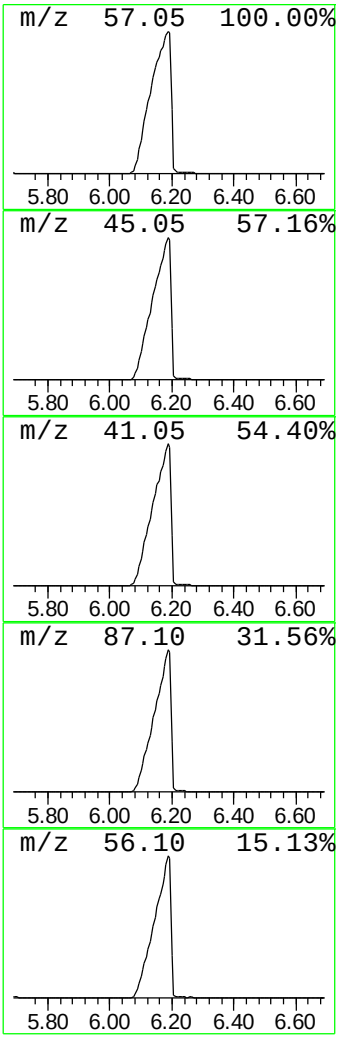
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.19	1637.26 ng	22895200	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
2	Formic acid, 3,3-dimethylbut-2-y...	130	C7H14O2	1000368-20-5	45
3	Propane, 1-methoxy-2,2-dimethyl-	102	C6H14O	001118-00-9	42
4	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	40
5	1,4-Pentanediol	104	C5H12O2	000626-95-9	38





Library Search Compound Report

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Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

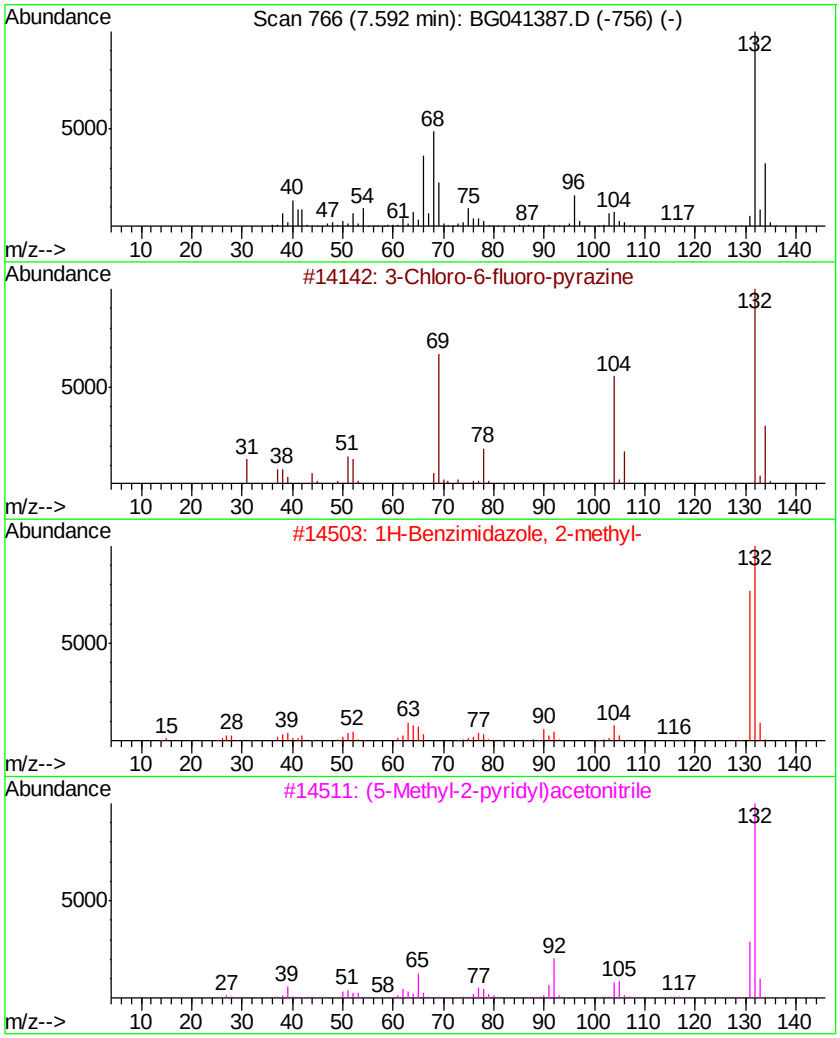
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

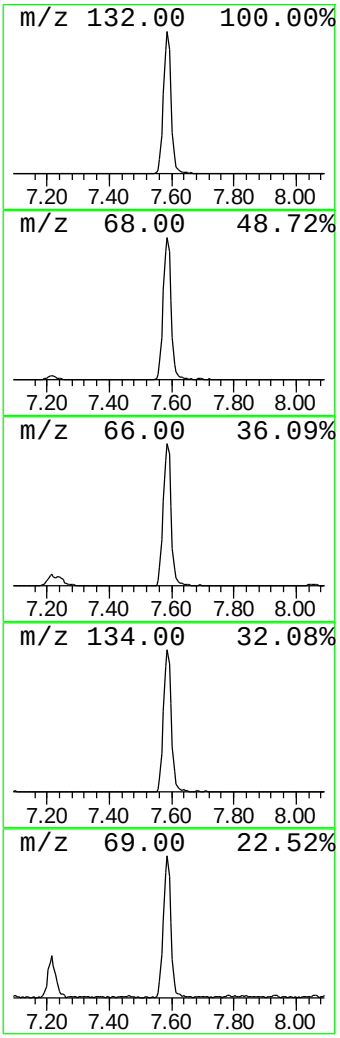
****

Peak Number 3 unknown7.59 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	64.66 ng	904260	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3	(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	22
5	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14





Library Search Compound Report

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Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

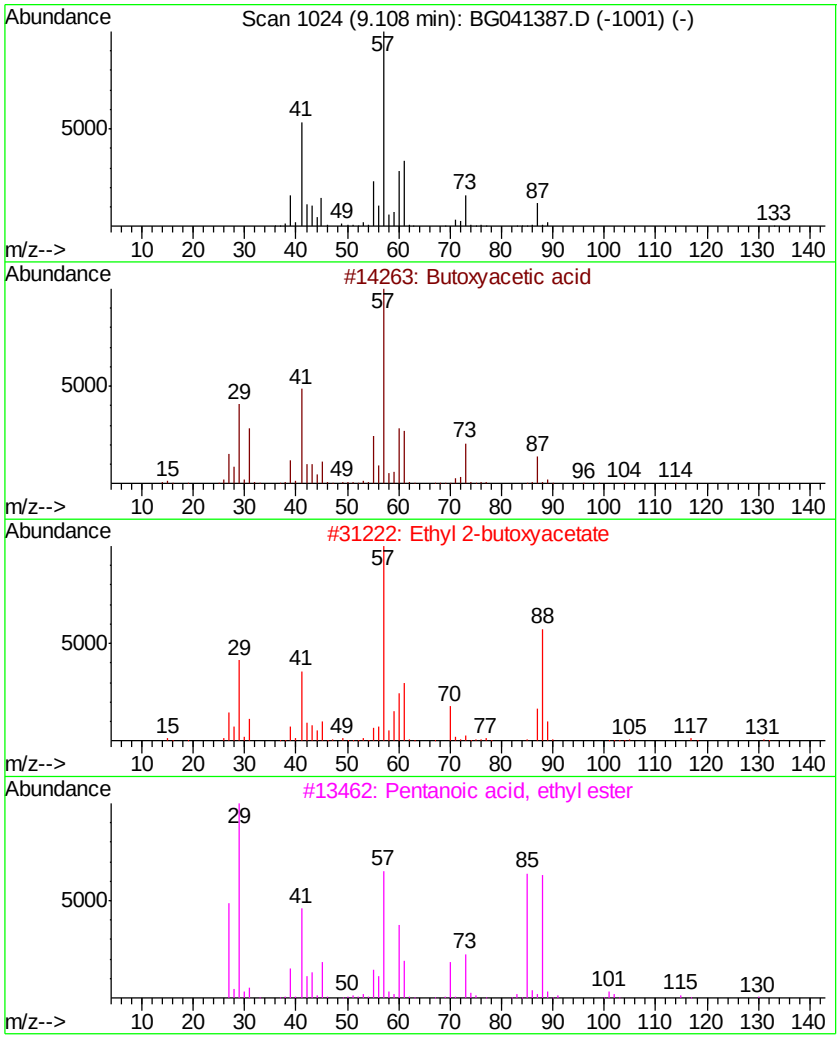
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

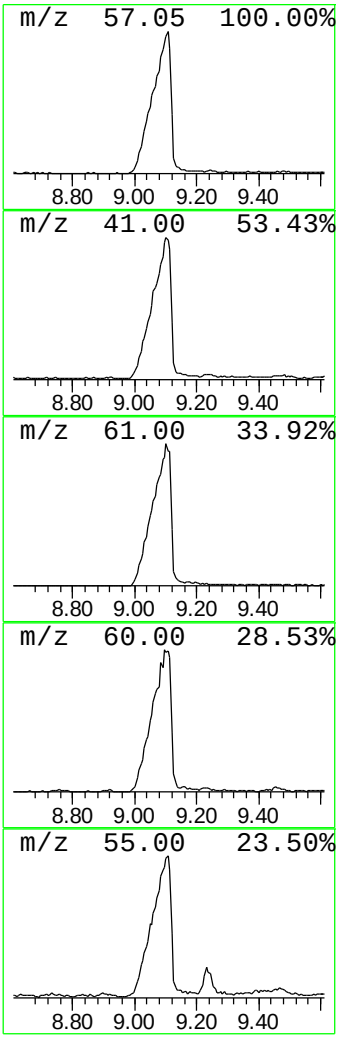
****

Peak Number 5 Butoxyacetic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.11	103.77 ng	1451130	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butoxyacetic acid	132	C6H12O3	002516-93-0	78
2	Ethyl 2-butoxyacetate	160	C8H16O3	1000366-89-3	42
3	Pentanoic acid, ethyl ester	130	C7H14O2	000539-82-2	40
4	Propanoic acid, 2,2-dimethyl-, s...	208	C5H9AgO2	007324-58-5	35
5	Butyl glycolate	132	C6H12O3	007397-62-8	28





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

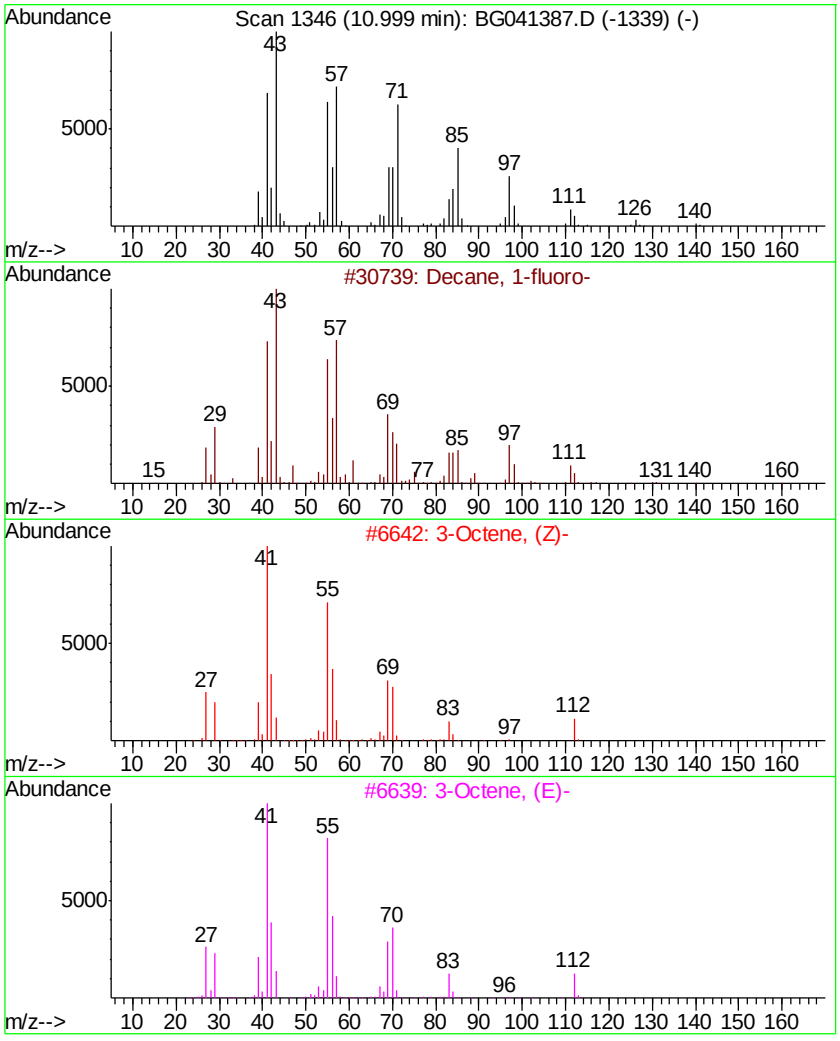
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

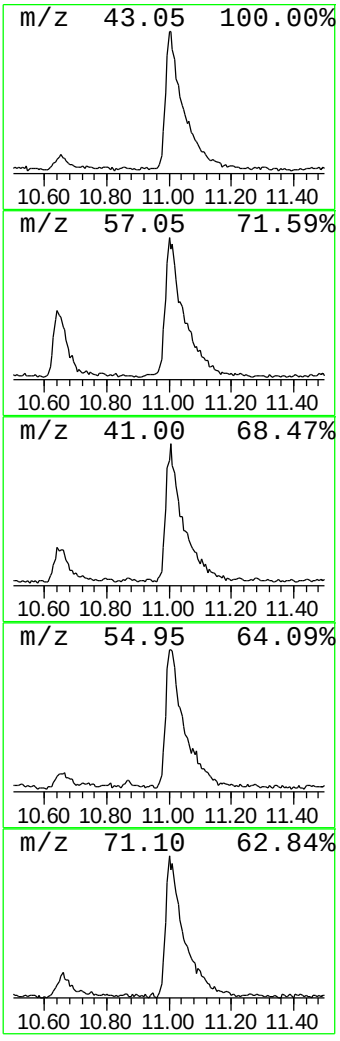
****

Peak Number 6 Decane, 1-fluoro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	42.24 ng	727994	Naphthalene-d8	10.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 1-fluoro-	160	C10H21F	000334-56-5	58
2	3-Octene, (Z)-	112	C8H16	014850-22-7	38
3	3-Octene, (E)-	112	C8H16	014919-01-8	38
4	2-Propenoic acid, octyl ester	184	C11H20O2	002499-59-4	35
5	1-Hexene, 4-ethyl-	112	C8H16	016746-85-3	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

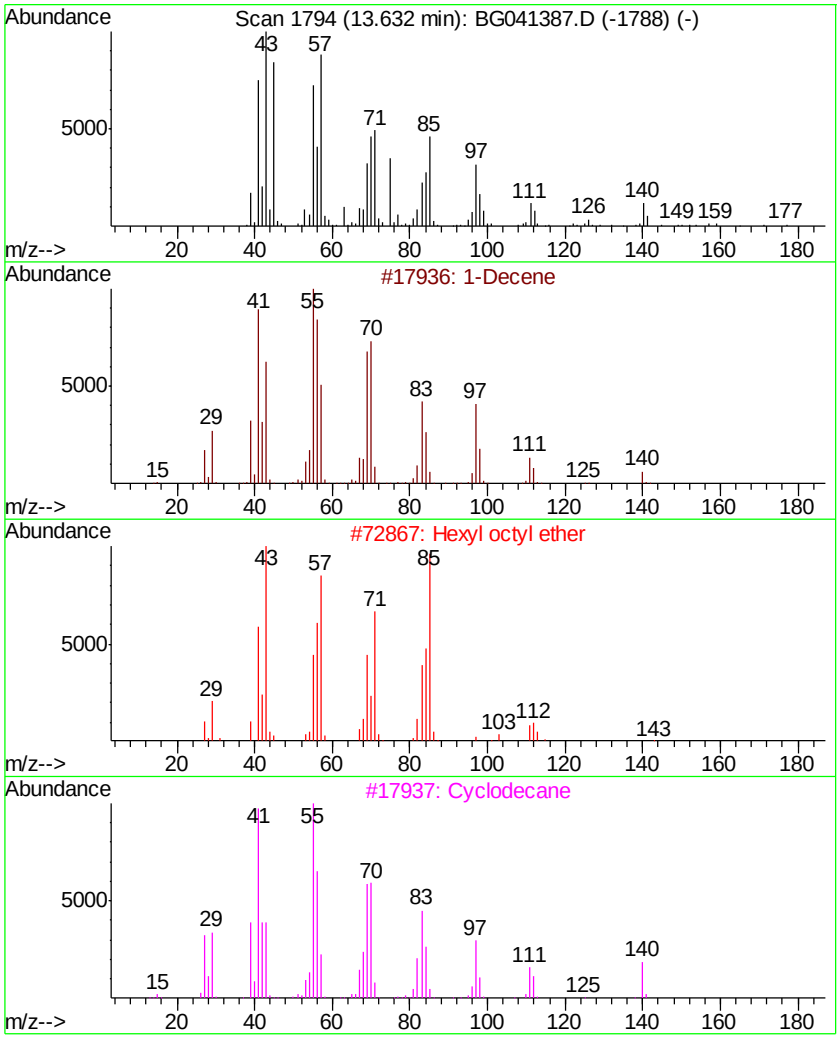
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

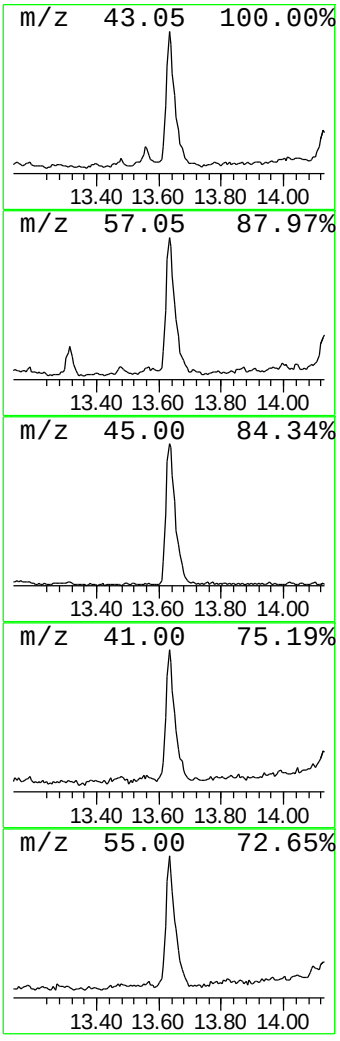
****

Peak Number 7 unknown13.63 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	23.09 ng	586885	Acenaphthene-d10	14.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decene	140	C10H20	000872-05-9	46
2	Hexyl octyl ether	214	C14H30O	017071-54-4	43
3	Cyclodecane	140	C10H20	000293-96-9	38
4	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	27
5	(2,2-Dimethylcyclobutyl)methylamine	113	C7H15N	1000195-04-0	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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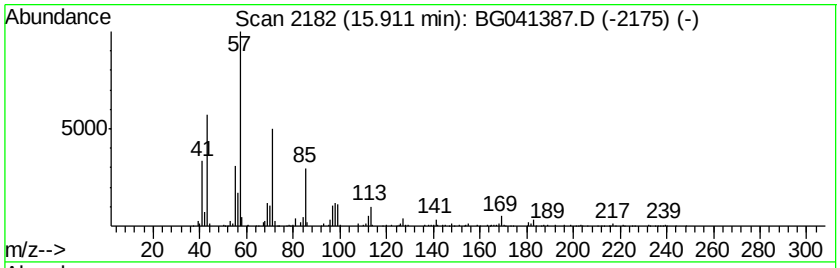
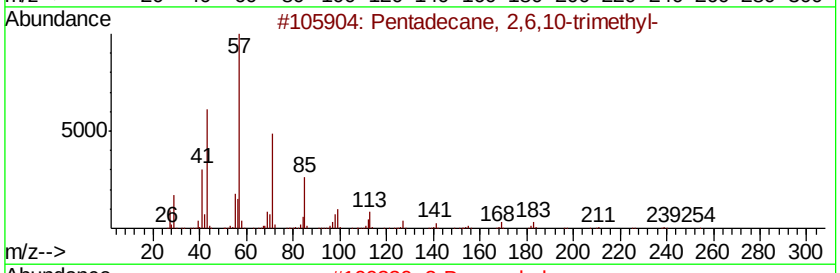
****

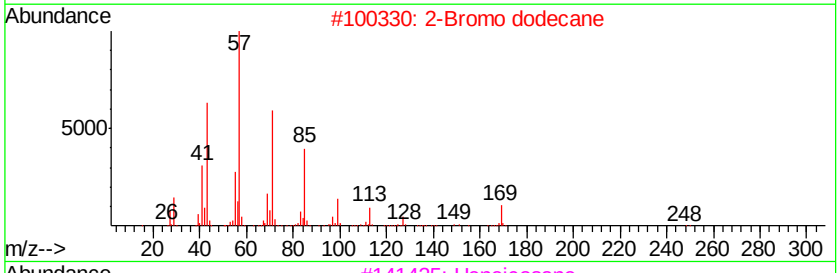
Peak Number 8 Pentadecane, 2,6,10-trimethyl- Concentration Rank 7

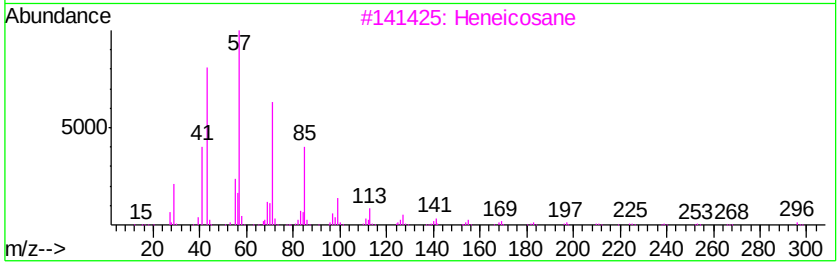
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	41.29 ng	1049360	Acenaphthene-d10	14.70

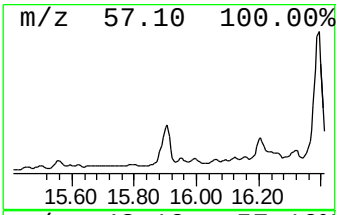
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	94
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	74
3	Heneicosane	296	C21H44	000629-94-7	72
4	Tridecane, 6-propyl-	226	C16H34	055045-10-8	70
5	3,5-Dimethyldodecane	198	C14H30	107770-99-0	70

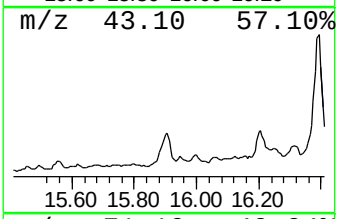


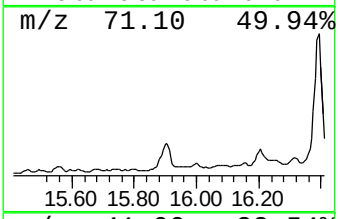


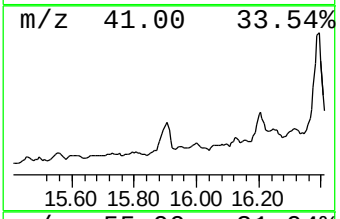


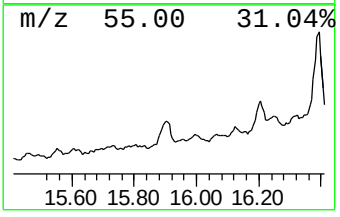












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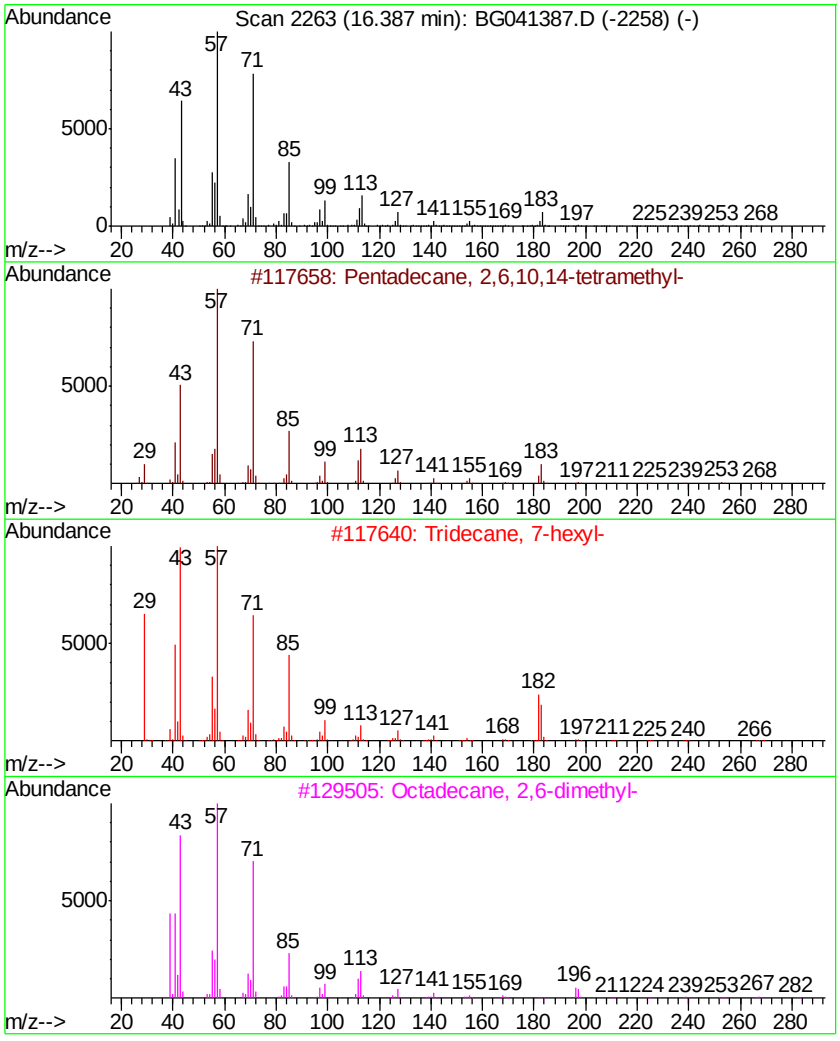
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

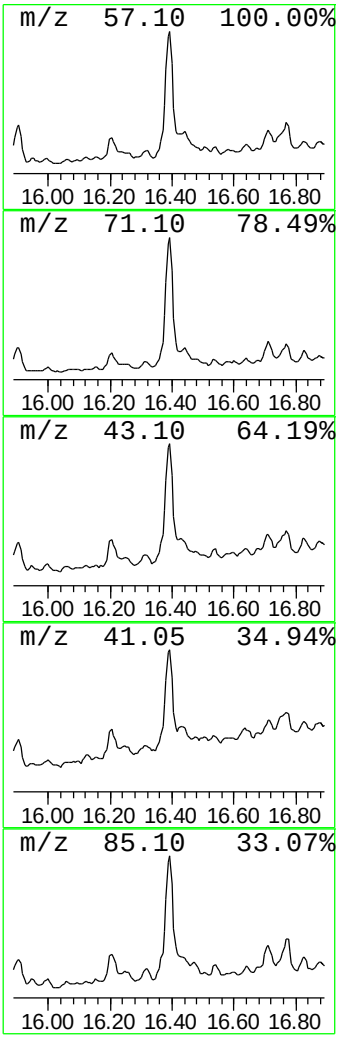
****

Peak Number 9 Pentadecane, 2,6,10,14-tetr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.39	19.48 ng	3850090	Phenanthrene-d10	17.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90
3	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	90
4	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	87
5	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

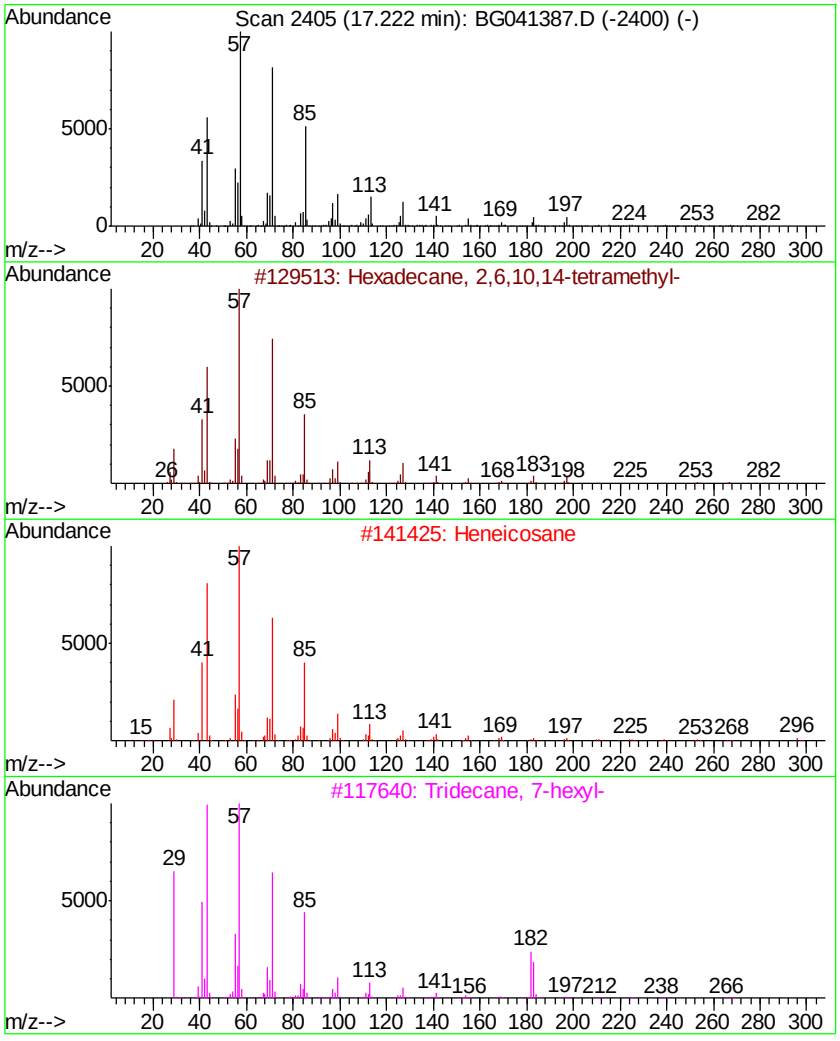
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

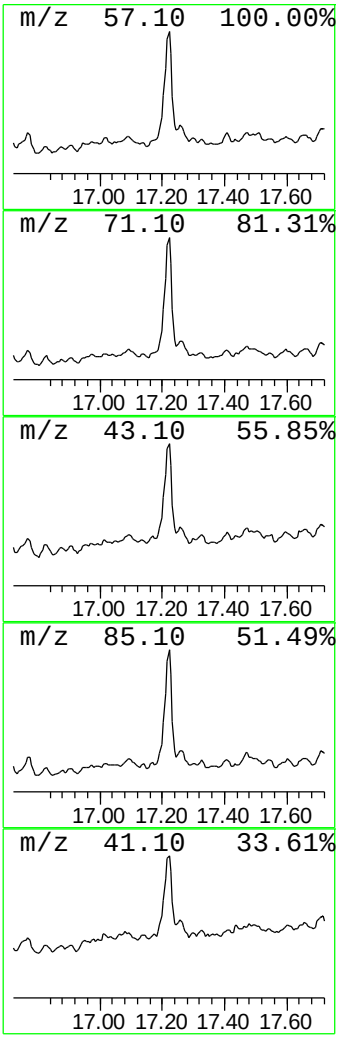
****

Peak Number 10 Hexadecane, 2,6,10,14-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.22	20.00 ng	3952000	Phenanthrene-d10	17.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	96
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87
4	Nonadecane	268	C19H40	000629-92-5	87
5	Hexacosane	366	C26H54	000630-01-3	86





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Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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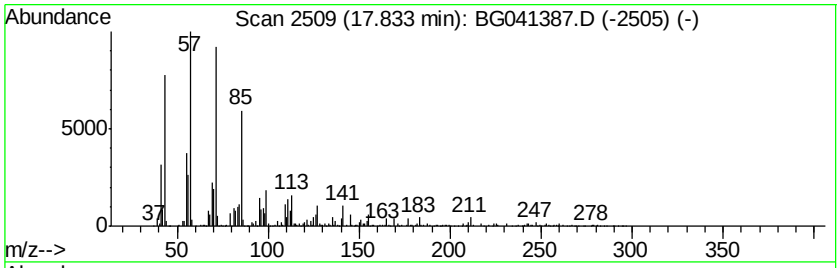
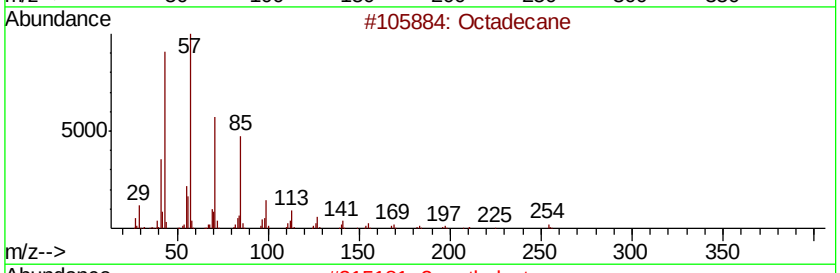
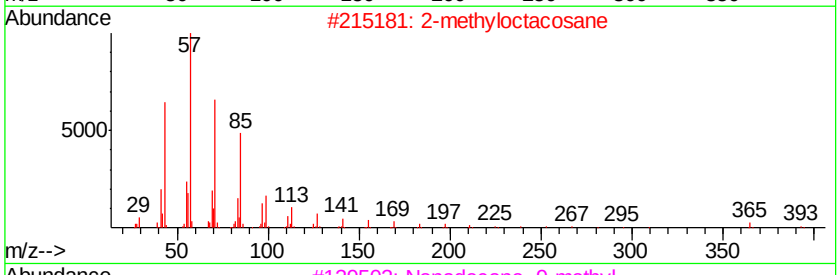
Peak Number 11 Octadecane Concentration Rank 16

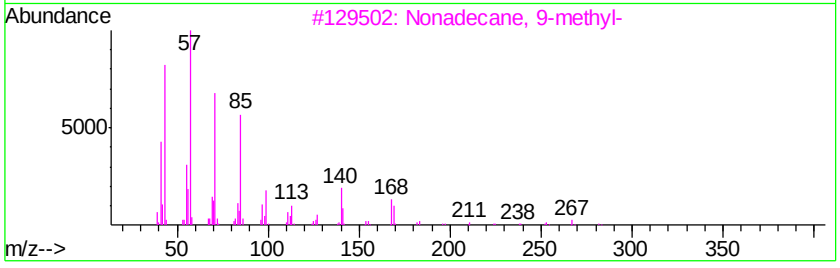
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	7.38 ng	1458560	Phenanthrene-d10	17.45

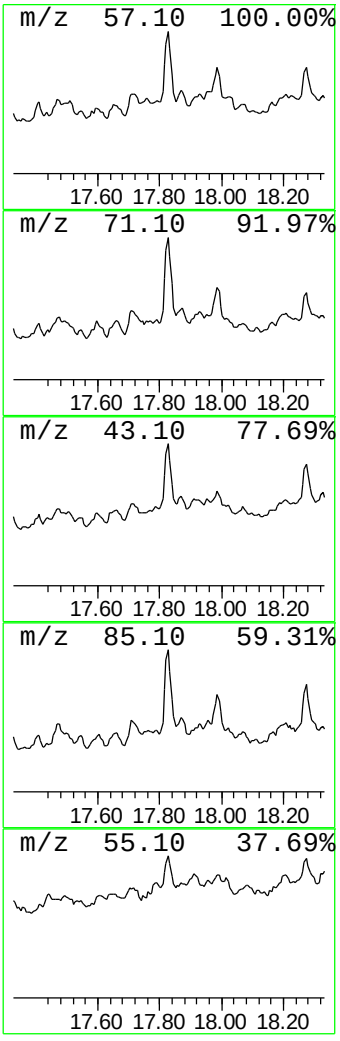
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
4	Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	91
5	Heneicosane	296	C21H44	000629-94-7	91











Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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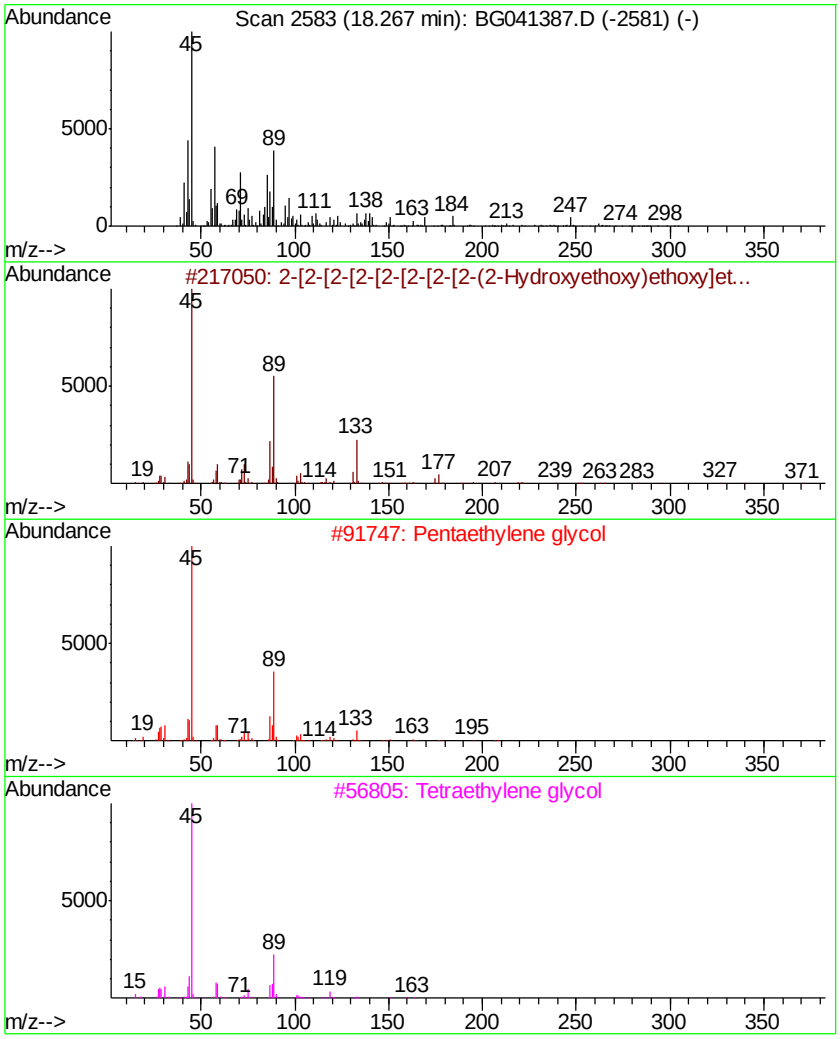
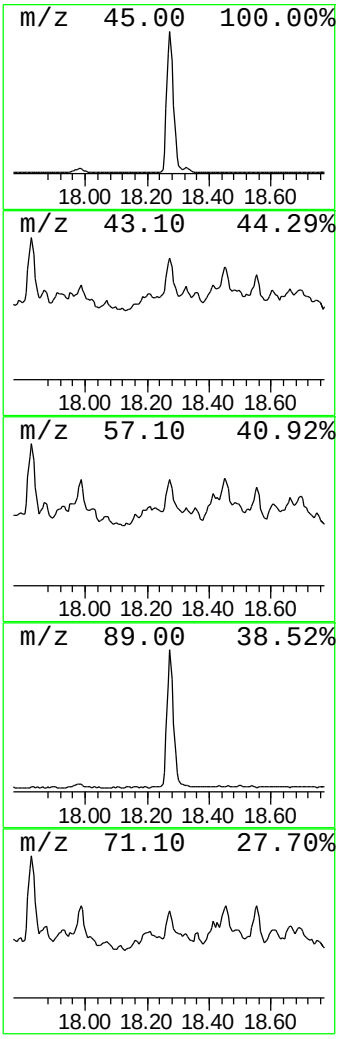
Peak Number 12 2-[2-[2-[2-[2-[2-[2-[2-(2-H... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.27	9.66 ng	1909030	Phenanthrene-d10	17.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-[2-[2-[2-[2-[2-[2-(2-Hydrox...	414	C18H38O10	1000352-12-5	58
2	Pentaethylene glycol	238	C10H22O6	004792-15-8	58
3	Tetraethylene glycol	194	C8H18O5	000112-60-7	53
4	Hexaethylene glycol	282	C12H26O7	002615-15-8	53
5	2-[2-[2-[2-[2-[2-[2-[2-[2-(2-Hyd...	458	C20H42O11	1000352-12-6	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

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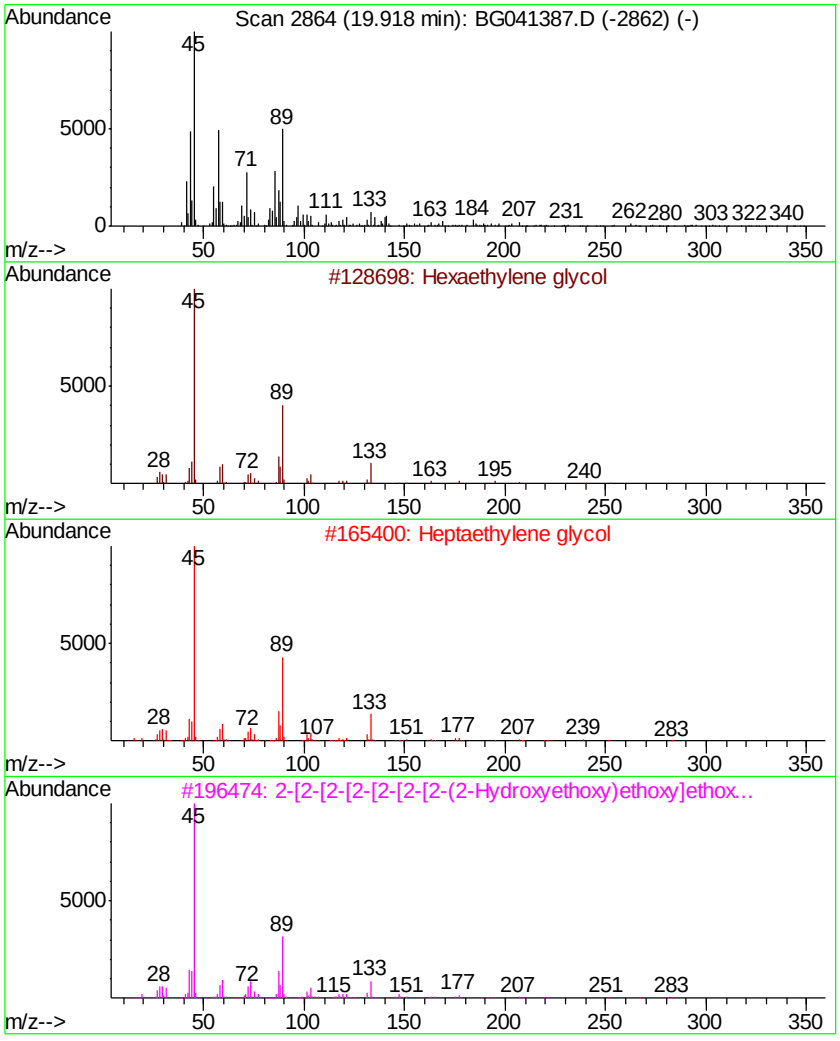
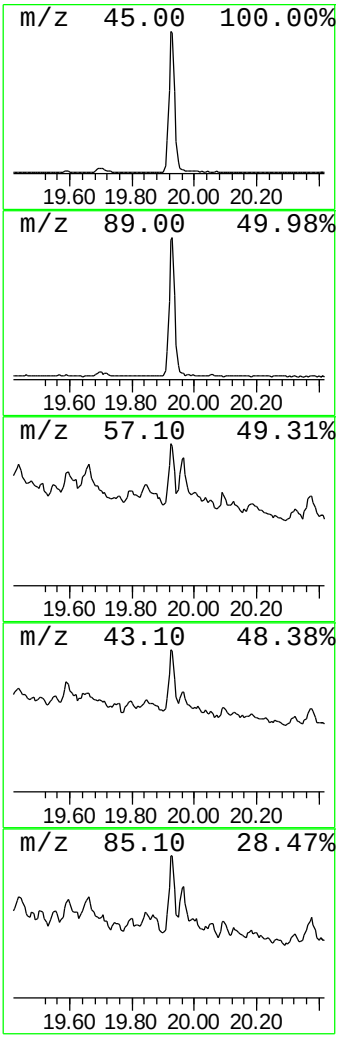
Peak Number 13 Hexaethylene glycol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.92	15.31 ng	1868920	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexaethylene glycol	282	C12H26O7	002615-15-8	68
2	Heptaethylene glycol	326	C14H30O8	005617-32-3	68
3	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	68
4	Pentaethylene glycol	238	C10H22O6	004792-15-8	68
5	2-[2-[2-[2-[2-[2-[2-(2-Hydrox...	414	C18H38O10	1000352-12-5	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062119\
Data File : BG041387.D
Acq On : 21 Jun 2019 21:51
Operator : HP/JU
Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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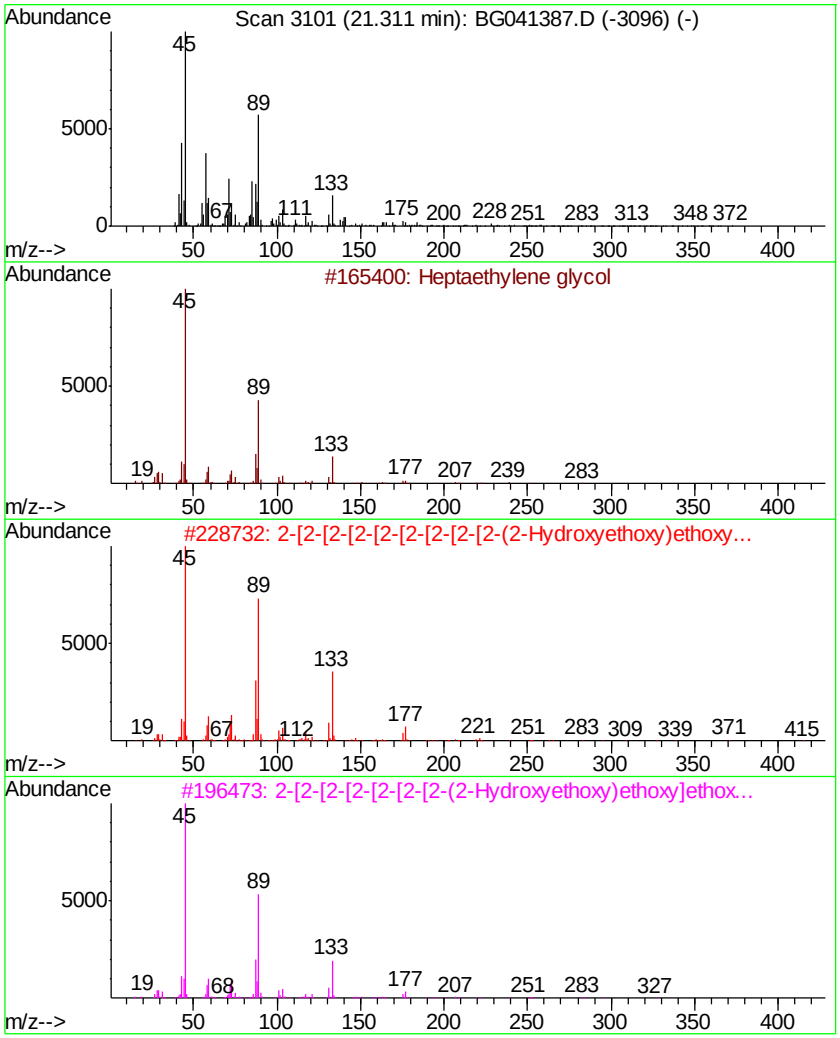
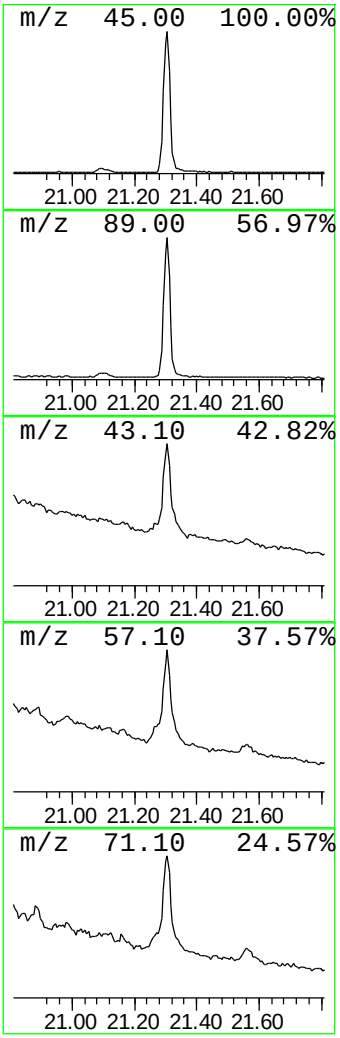
Peak Number 14 Heptaethylene glycol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.31	20.00 ng	2440820	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptaethylene glycol	326	C14H30O8	005617-32-3	76
2	2-[2-[2-[2-[2-[2-[2-[2-(2-Hyd...	458	C20H42O11	1000352-12-6	74
3	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	74
4	2-[2-[2-[2-[2-[2-[2-[2-[2-[2-...	546	C24H50O13	1000352-12-7	74
5	2-[2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	68





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Sample : K3433-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061719.M
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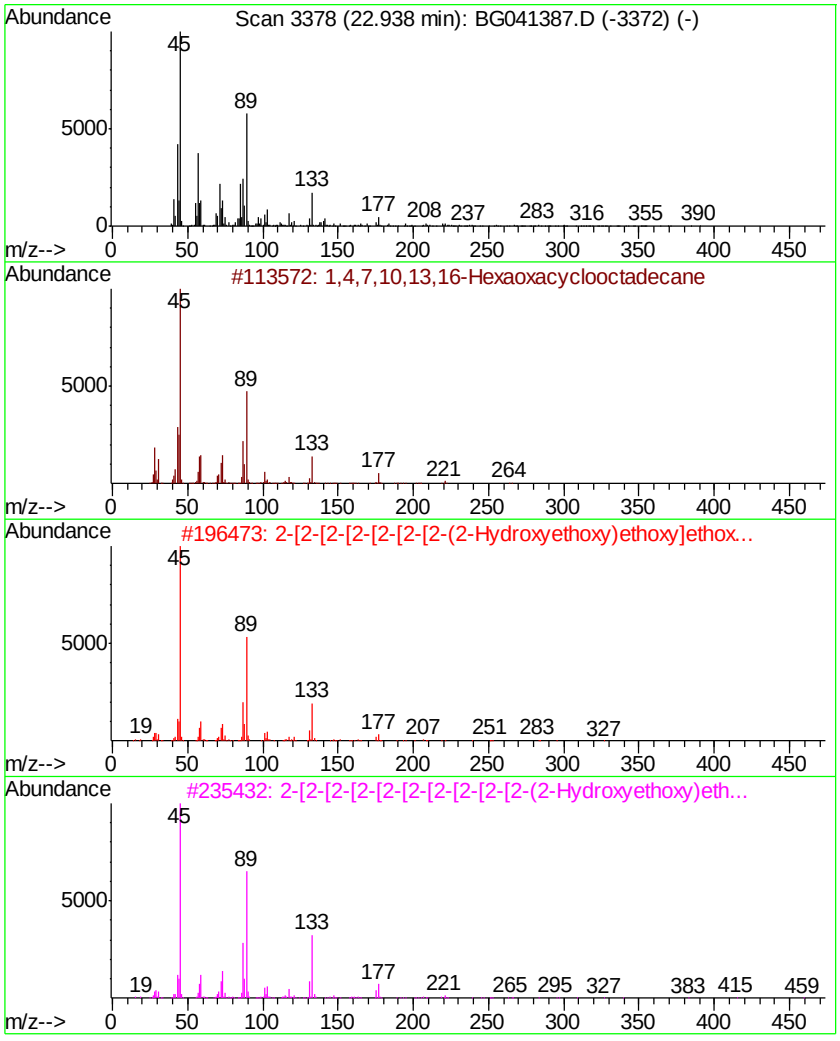
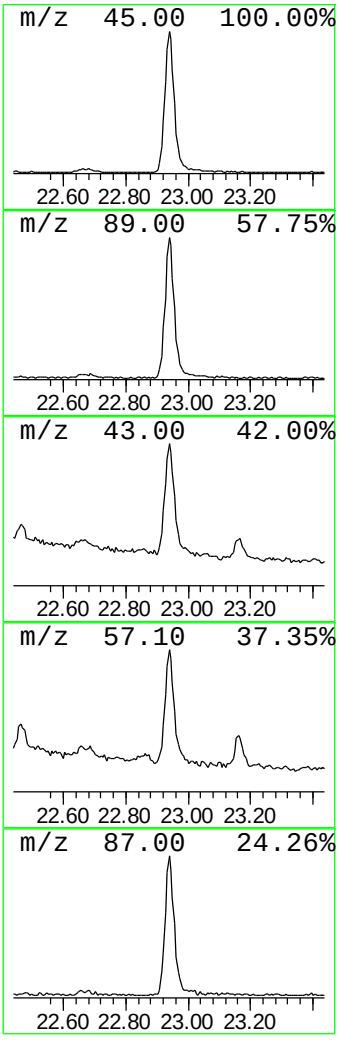
Peak Number 15 1,4,7,10,13,16-Hexaoxacyclo... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.94	11.43 ng	1395100	Chrysene-d12	21.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4,7,10,13,16-Hexaoxacyclooctad...	264	C12H24O6	017455-13-9	93
2	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	81
3	2-[2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	76
4	2-[2-[2-[2-[2-[2-[2-[2-(2-Hyd...	458	C20H42O11	1000352-12-6	74
5	Hexaethylene glycol	282	C12H26O7	002615-15-8	74





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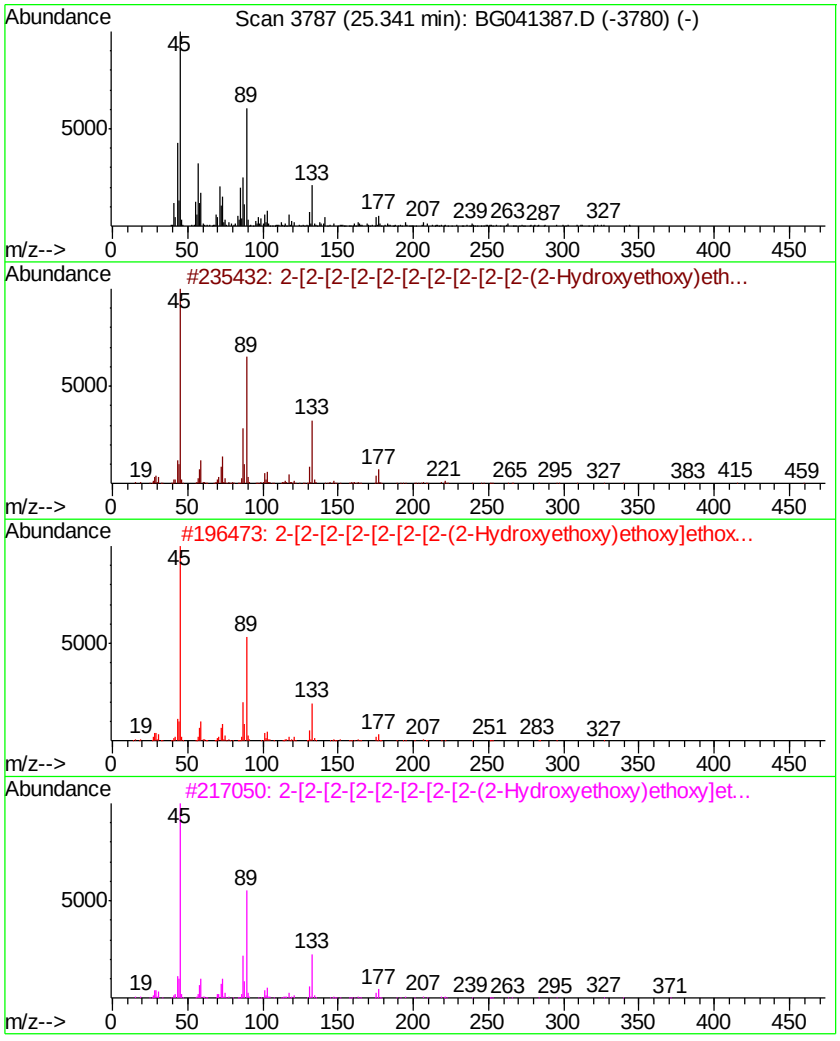
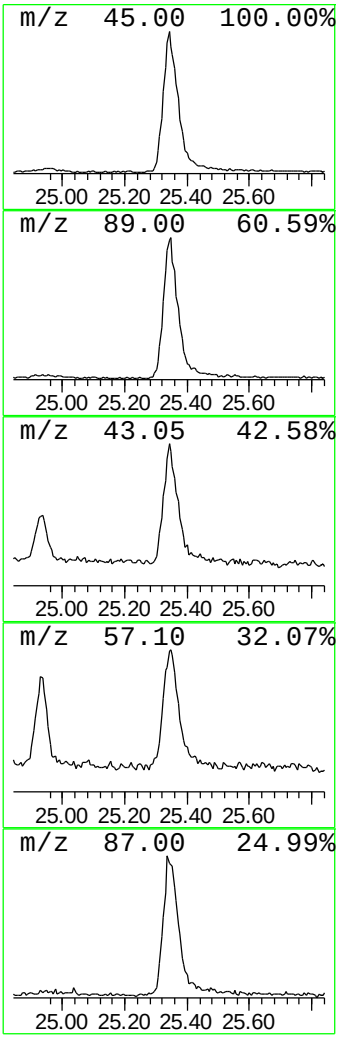
Peak Number 16 2-[2-[2-[2-[2-[2-[2-[2-[2-[... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.34	29.56 ng	923391	Perylene-d12	25.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-[2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	81
2	2-[2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	81
3	2-[2-[2-[2-[2-[2-[2-(2-Hydrox...	414	C18H38O10	1000352-12-5	81
4	Heptaethylene glycol	326	C14H30O8	005617-32-3	74
5	2-[2-[2-[2-[2-[2-[2-[2-[2-[2-...	546	C24H50O13	1000352-12-7	72





Tentatively Identified Compound (LSC) summary

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TIC Library : C:\Database\NIST11.L
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					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethanol, 2-butoxy-	6.19	1637.3	ng	22895200	1	8.06	279676	20.0
unknown7.59	7.59	64.7	ng	904260	1	8.06	279676	20.0
Butoxyacetic acid	9.11	103.8	ng	1451130	1	8.06	279676	20.0
Decane, 1-fluoro-	11.00	42.2	ng	727994	2	10.88	344694	20.0
unknown13.63	13.63	23.1	ng	586885	3	14.70	508322	20.0
Pentadecane, 2,6,...	15.91	41.3	ng	1049360	3	14.70	508322	20.0
Pentadecane, 2,6,...	16.39	19.5	ng	3850090	4	17.45	3952000	20.0
Hexadecane, 2,6,1...	17.22	20.0	ng	3952000	4	17.45	3952000	20.0
Octadecane	17.83	7.4	ng	1458560	4	17.45	3952000	20.0
2-[2-[2-[2-[2-[2-...	18.27	9.7	ng	1909030	4	17.45	3952000	20.0
Hexaethylene glycol	19.92	15.3	ng	1868920	5	21.73	2440820	20.0
Heptaethylene glycol	21.31	20.0	ng	2440820	5	21.73	2440820	20.0
1,4,7,10,13,16-He...	22.94	11.4	ng	1395100	5	21.73	2440820	20.0
2-[2-[2-[2-[2-[2-...	25.34	29.6	ng	923391	6	25.02	624816	20.0