LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

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_P\METHODS\8270E-BP110320.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004046.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	25	32	42	rBV	188505	364439	6.30%	1.208%
2	3.164	42	47	62	rVB	881817	1187534	20.54%	3.936%
3	3.840	157	162	171	rVB	58525	85753	1.48%	0.284%
4	5.264	399	404	411	rBV	101205	142915	2.47%	0.474%
5	5.805	488	496	506	rBV	1018083	1601096	27.69%	5.306%
6	6.399	592	597	606	rVB	50157	75969	1.31%	0.252%
7	7.452	769	776	791	rBV	589390	1035844	17.91%	3.433%
8	7.840	836	842	848	rBV	38443	62920	1.09%	0.209%
9	8.263	907	914	923	rVB	408853	641927	11.10%	2.127%
10	9.428	1104	1112	1133	rBV	1360496	2272826	39.31%	7.532%
11	9.946	1196	1200	1210	rVB	38689	64730	1.12%	0.215%
12	11.093	1387	1395	1402	rBV	502371	847932	14.66%	2.810%
13	13.510	1798	1806	1816	rBV	3231594	4551385	78.71%	15.084%
14	13.593	1816	1820	1829	rVB2	32492	64105	1.11%	0.212%
15	13.775	1845	1851	1864	rVB	323111	512806	8.87%	1.700%
16	14.310	1937	1942	1947	rBV	67031	91483	1.58%	0.303%
17	14.887	2033	2040	2048	rBV	721674	1063817	18.40%	3.526%
18	15.986	2221	2227	2242	rBV	188434	324327	5.61%	1.075%
19	16.369	2286	2292	2313	rBV	2183253	3349457	57.93%	11.101%
20	17.616	2498	2504	2510	rBV	829422	1223557	21.16%	4.055%
21	17.898	2548	2552	2562	rBV2	41023	70424	1.22%	0.233%
22	18.463	2644	2648	2666	rBV	333953	717027	12.40%	2.376%
23	18.892	2717	2721	2726	rBV2	64596	86426	1.49%	0.286%
24	19.275	2782	2786	2793	rBV	164260	303220	5.24%	1.005%
25	19.569	2832	2836	2846	rVB2	133365	173215	3.00%	0.574%
26	19.669	2850	2853	2861	rBV	113556	184744	3.20%	0.612%
27	19.880	2885	2889	2892	rBV	47276	61131	1.06%	0.203%
28	20.116	2924	2929	2934	rBV	4865063	5782219	100.00%	19.163%
29	20.380	2968	2974	2981	rBV2	124670	266586	4.61%	0.884%
30	21.310	3128	3132	3141	rBV	245109	469525	8.12%	1.556%
31	21.651	3185	3190	3195	rBV	956663	1233428	21.33%	4.088%
32	24.121	3603	3610	3617	rVB	622225	1260965	21.81%	4.179%

Sum of corrected areas: 30173732

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
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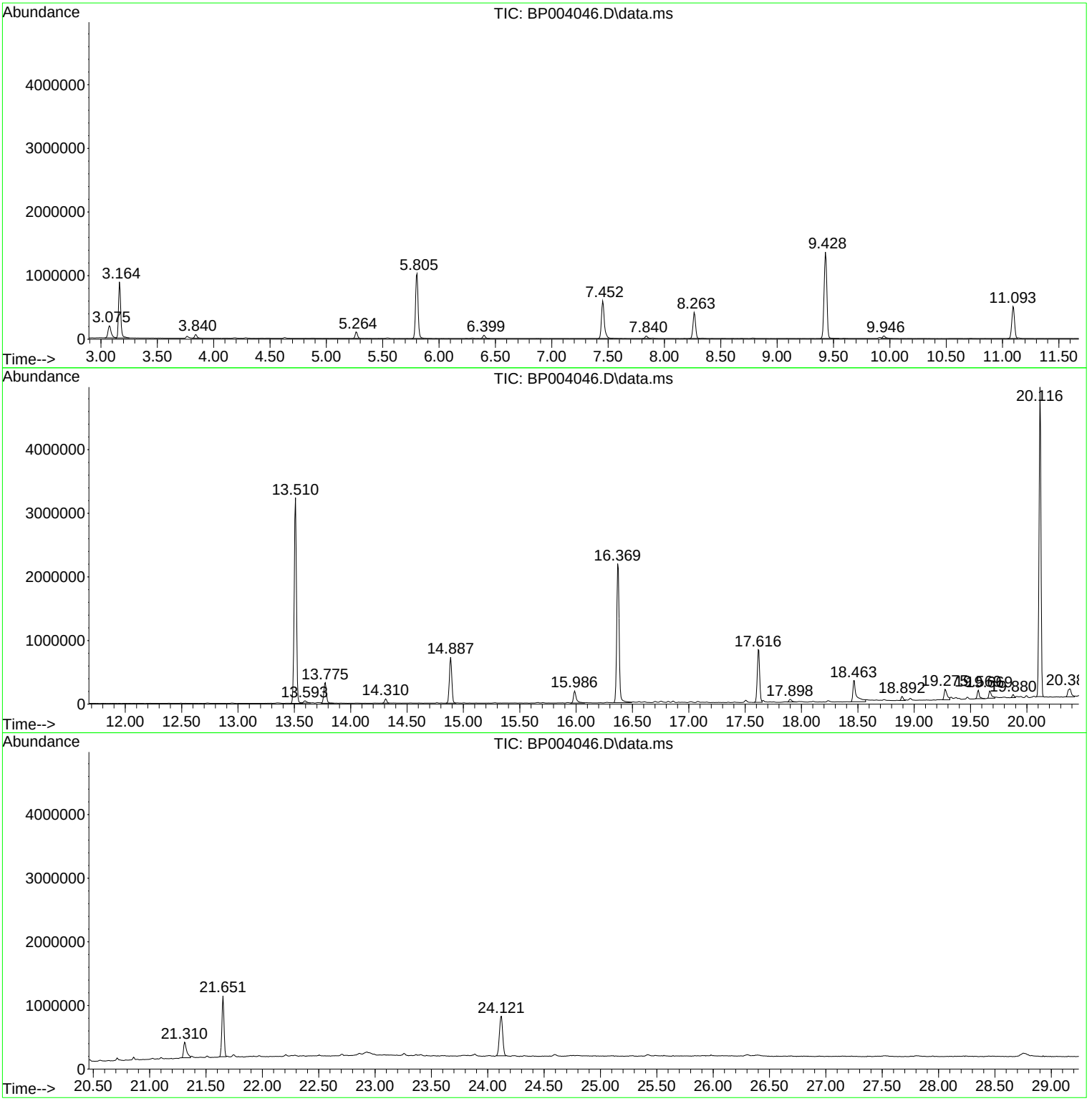
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

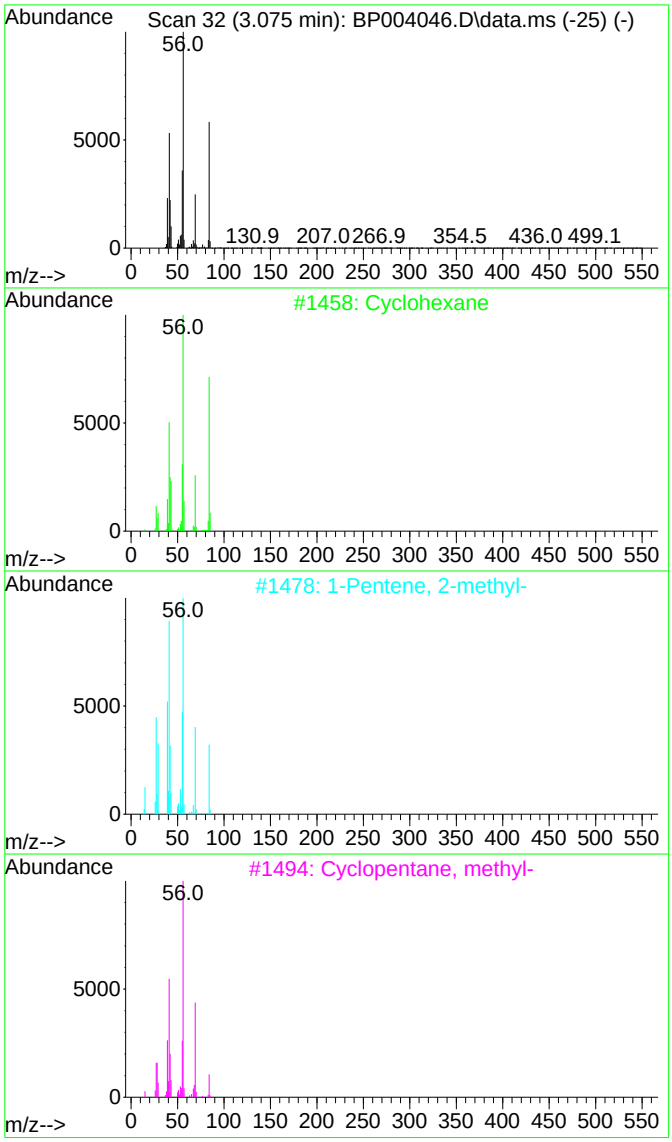
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

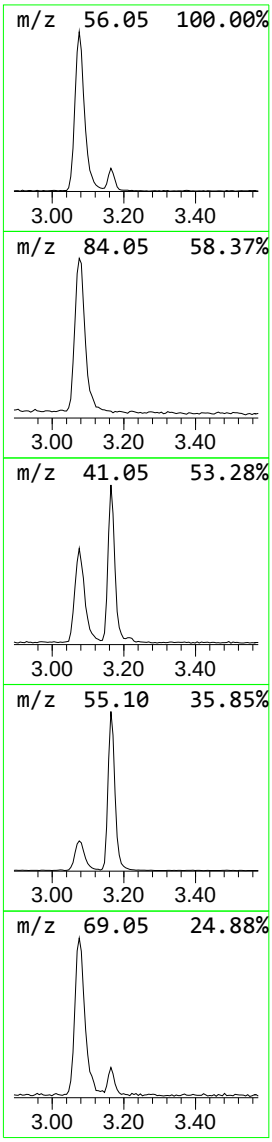
****

Peak Number 1 Cyclohexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	11.35 ng	364439	1,4-Dichlorobenzene-d4	8.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	91
2	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	59
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	47
5	1-Hexene, 5-methyl-	98	C7H14	003524-73-0	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

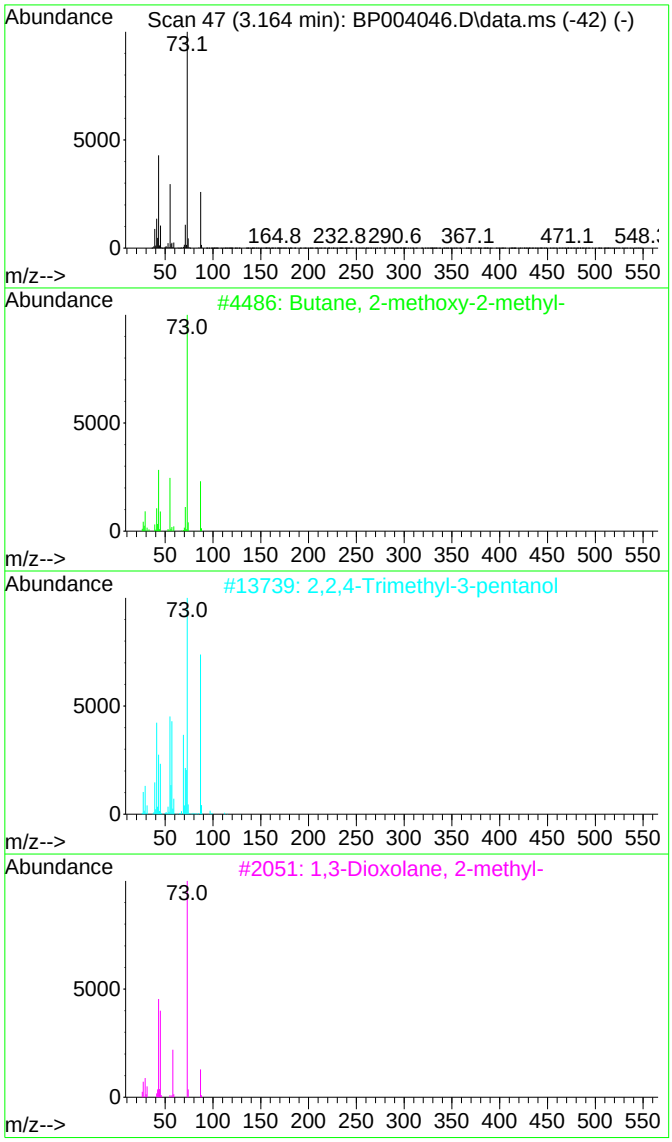
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

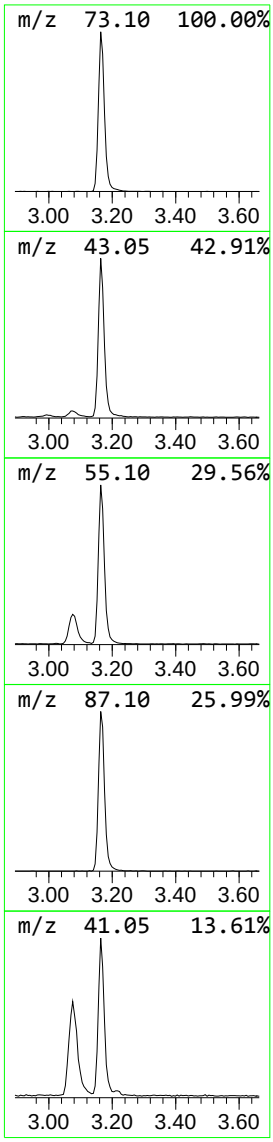
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.164	37.00 ng	1187530	1,4-Dichlorobenzene-d4	8.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	10





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Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

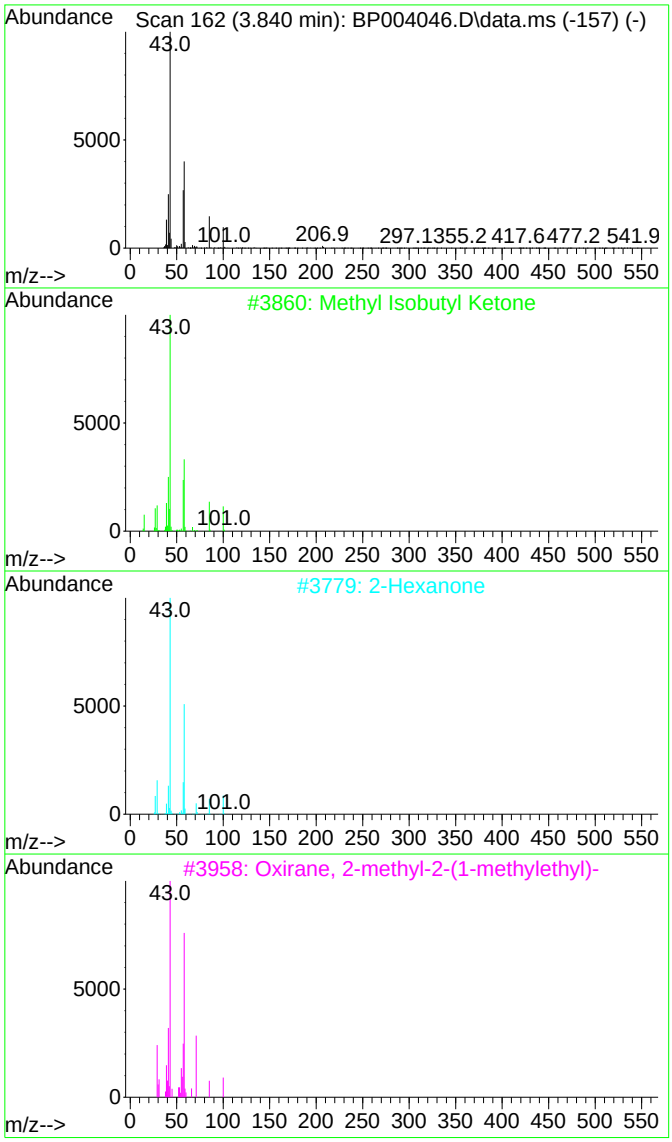
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

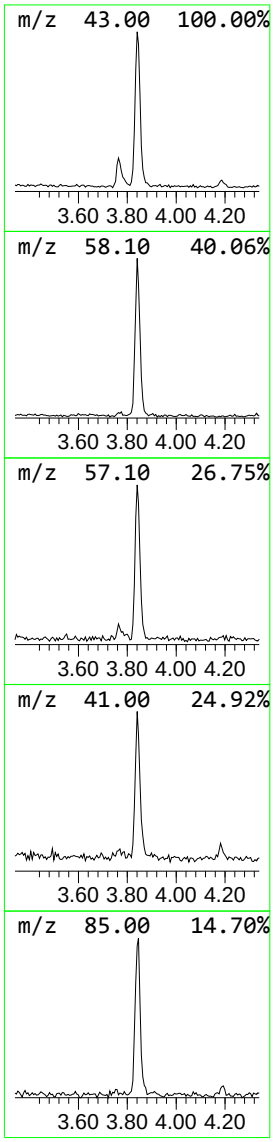
****

Peak Number 3 Methyl Isobutyl Ketone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.840	2.67 ng	85753	1,4-Dichlorobenzene-d4	8.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	86
2	2-Hexanone	100	C6H12O	000591-78-6	78
3	Oxirane, 2-methyl-2-(1-methyleth...	100	C6H12O	072221-03-5	64
4	Acetone	58	C3H6O	000067-64-1	40
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	37





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Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B5-GW

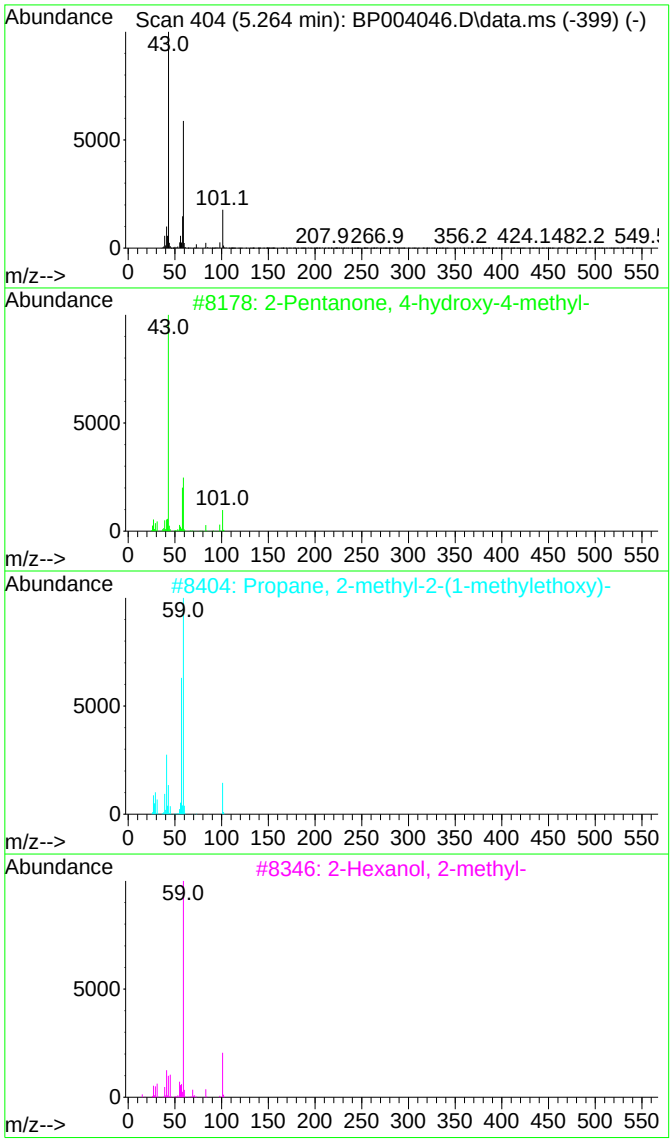
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

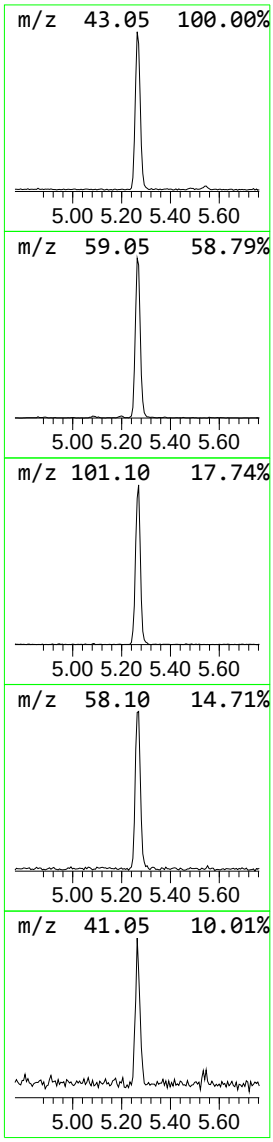
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.264	4.45 ng	142915	1,4-Dichlorobenzene-d4	8.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4	Tetraglyme	222	C10H22O5	000143-24-8	28
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
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Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B5-GW

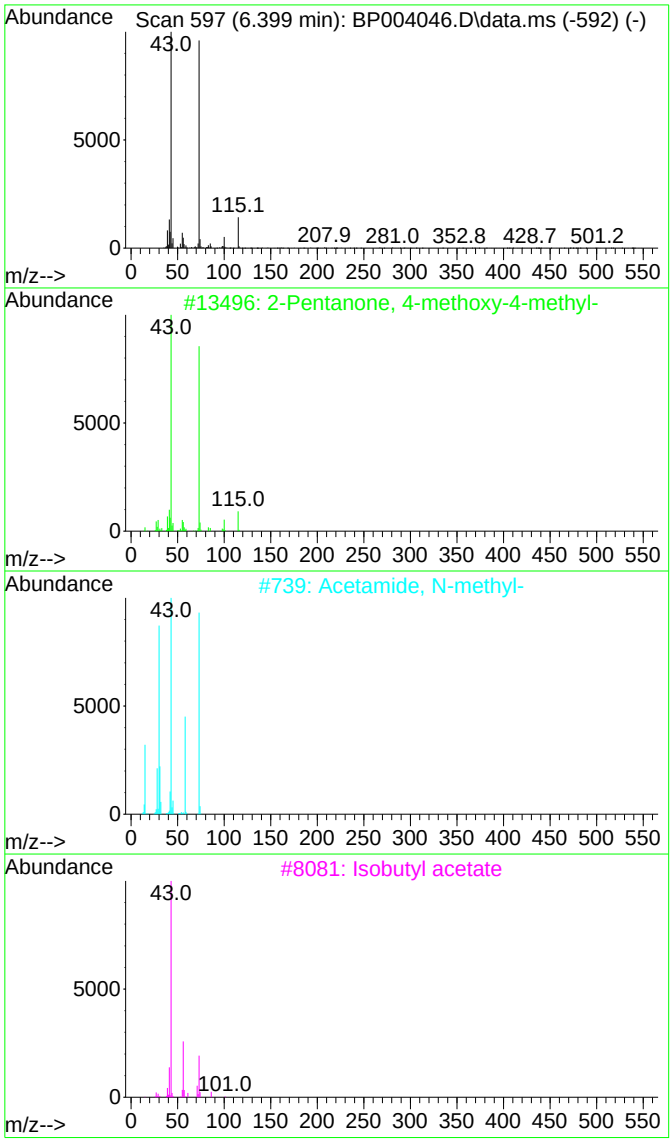
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

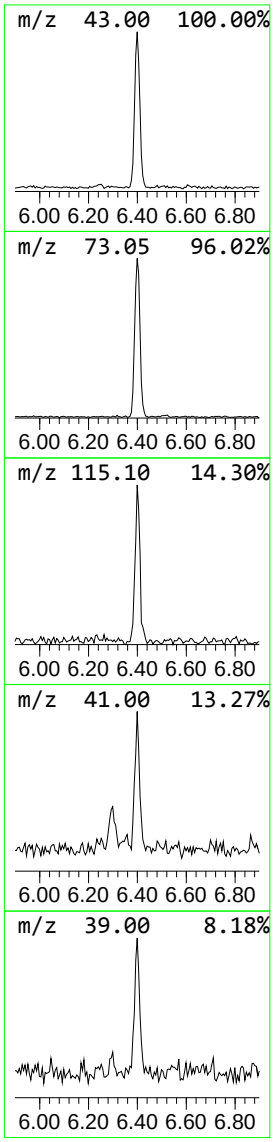
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown6.399 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.399	2.37 ng	75969	1,4-Dichlorobenzene-d4	8.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
2	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	47
3	Isobutyl acetate	116	C6H12O2	000110-19-0	23
4	D-chiro-Inositol, 3-O-(2-amino-4...	379	C14H25N3O9	006980-18-3	12
5	3-Methoxy-3-methylbutanol	118	C6H14O2	056539-66-3	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
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Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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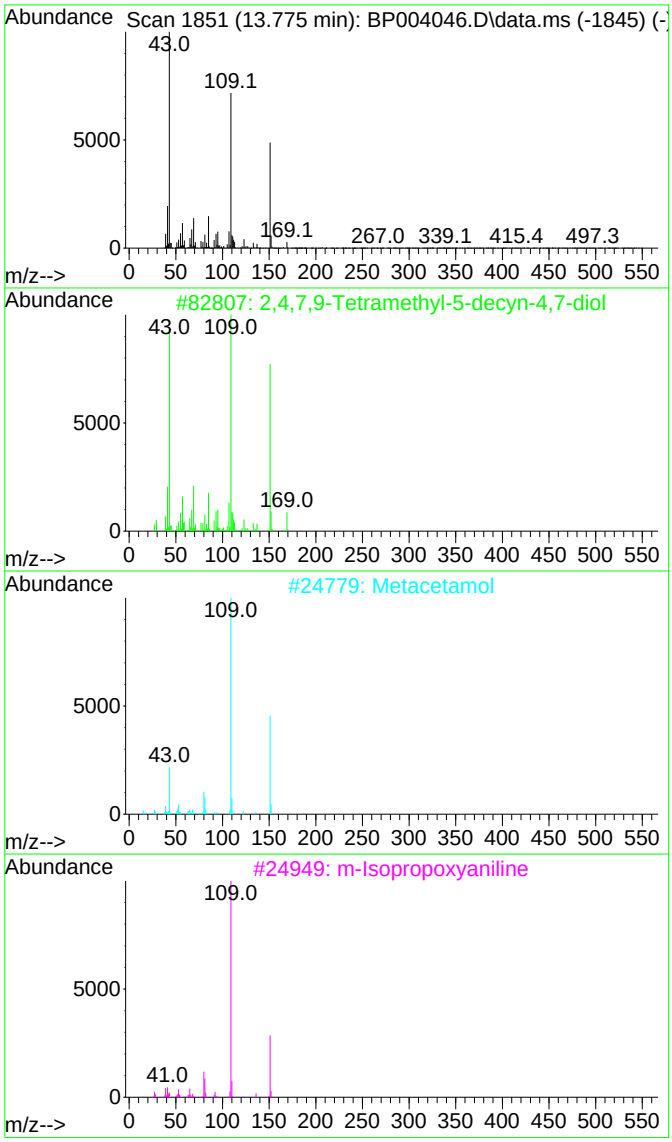
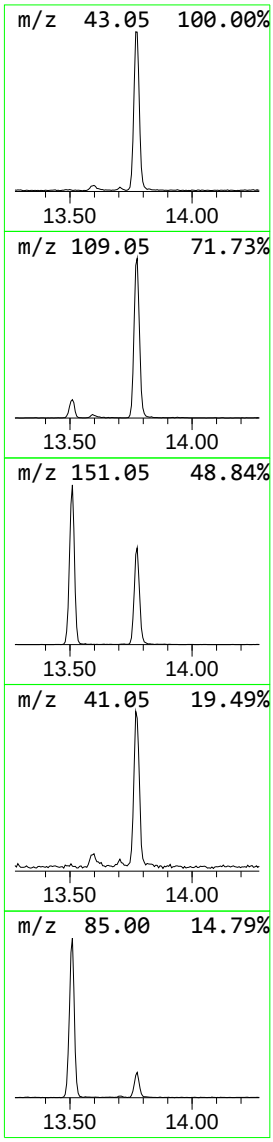
****
Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.775	9.64 ng	512806	Acenaphthene-d10	14.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	Metacetamol	151	C8H9NO2	000621-42-1	53
3	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	53
4	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53





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Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B5-GW

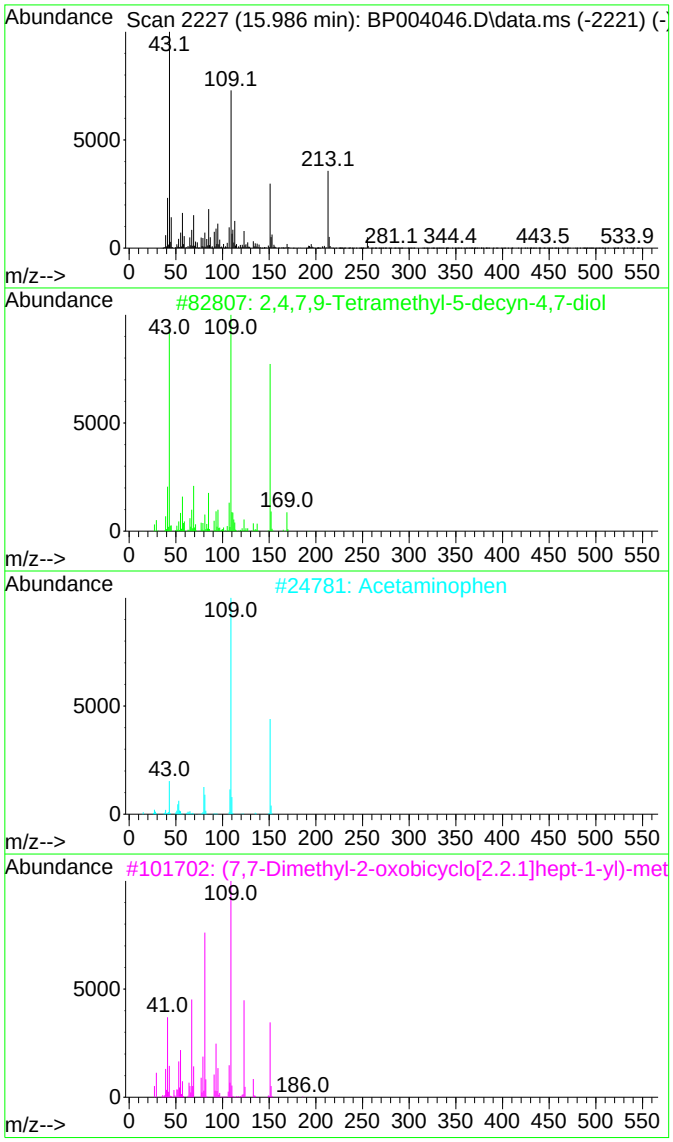
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

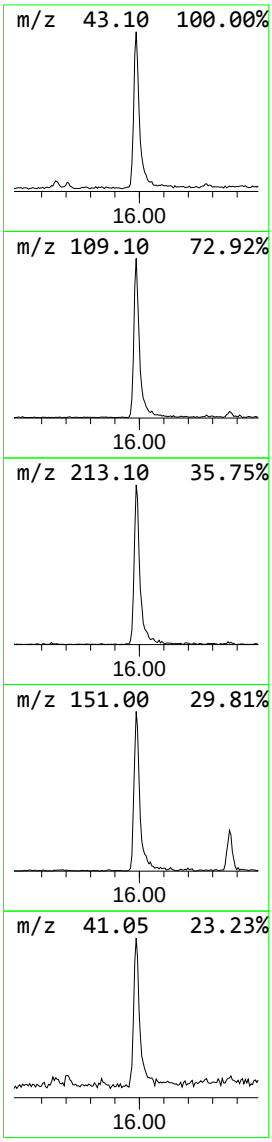
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 unknown15.986 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.986	6.10 ng	324327	Acenaphthene-d10	14.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	53
2	Acetaminophen	151	C8H9NO2	000103-90-2	38
3	(7,7-Dimethyl-2-oxobicyclo[2.2.1...	250	C10H15ClO3S	1000194-76-1	38
4	Cyclopentanol, 1-(methylenecyclo...	138	C9H14O	086951-58-8	38
5	Metacetamol	151	C8H9NO2	000621-42-1	35





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ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId : B5-GW

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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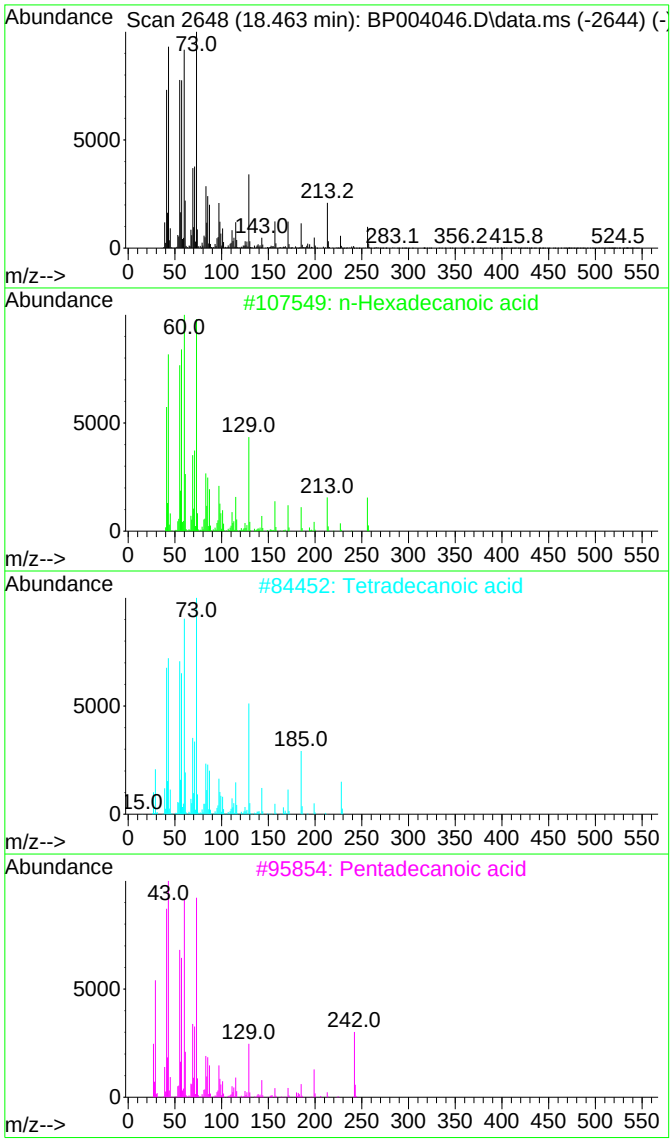
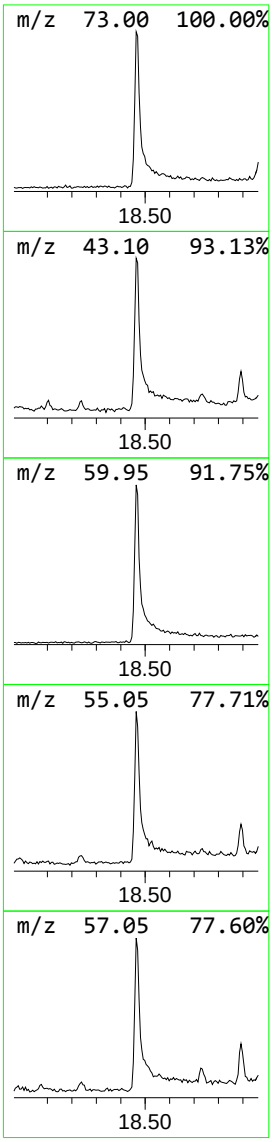
Peak Number 8 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.463	11.72 ng	717027	Phenanthrene-d10	17.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	Tridecanoic acid	214	C13H26O2	000638-53-9	81
5	Undecanoic acid	186	C11H22O2	000112-37-8	59





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Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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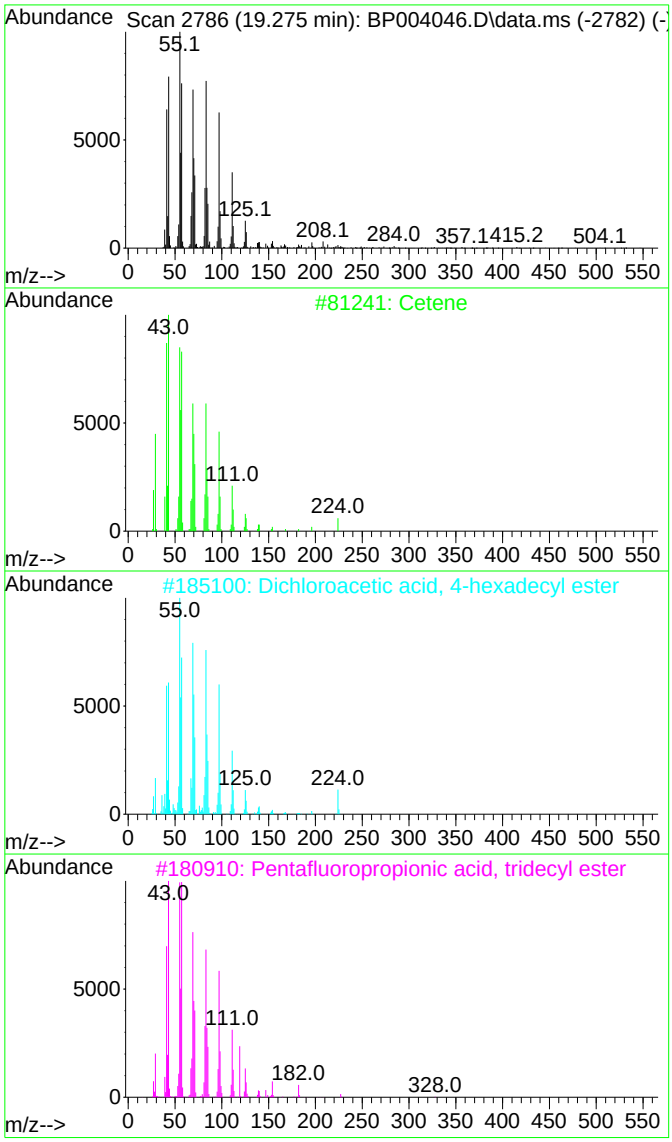
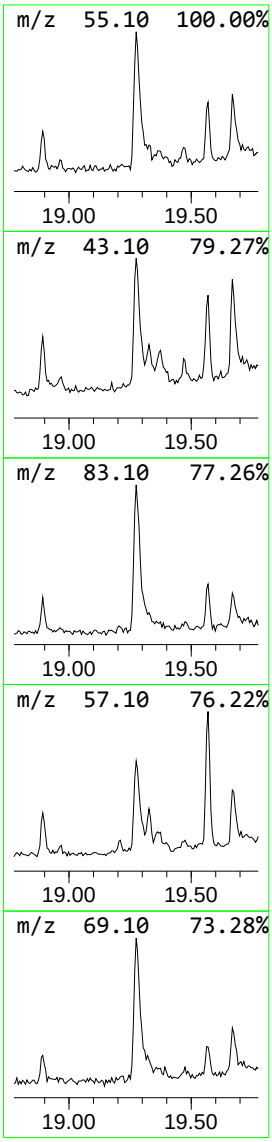
Peak Number 9 Cetene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.275	4.96 ng	303220	Phenanthrene-d10	17.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cetene	224	C16H32	000629-73-2	96
2	Dichloroacetic acid, 4-hexadecyl...	352	C18H34Cl2O2	1000282-98-1	94
3	Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	94
4	1-Tetradecanol	214	C14H30O	000112-72-1	91
5	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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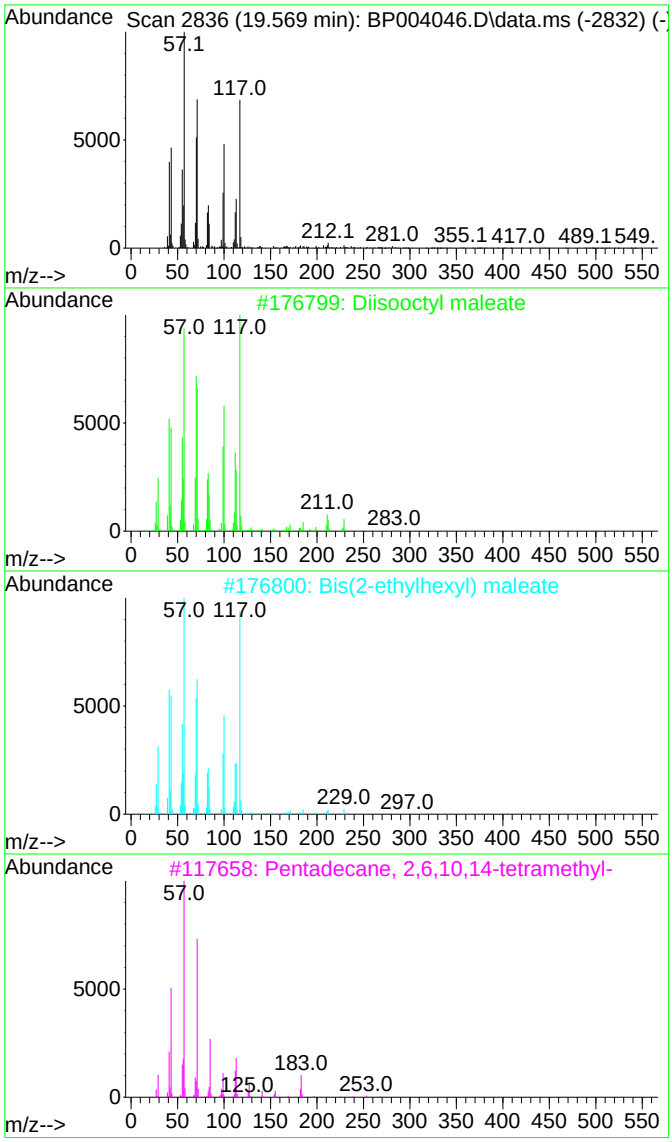
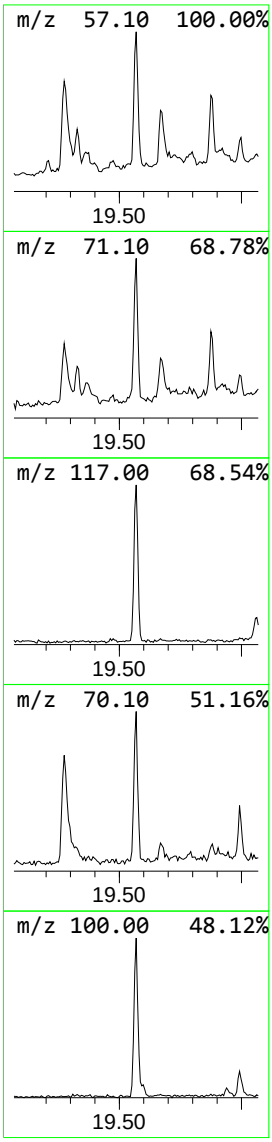
Peak Number 10 Diisooctyl maleate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.569	2.83 ng	173215	Phenanthrene-d10	17.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisooctyl maleate	340	C20H36O4	001330-76-3	83
2	Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	64
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	42
4	Oxalic acid, 2-ethylhexyl ethyl ...	230	C12H22O4	1000309-38-4	38
5	Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000357-82-9	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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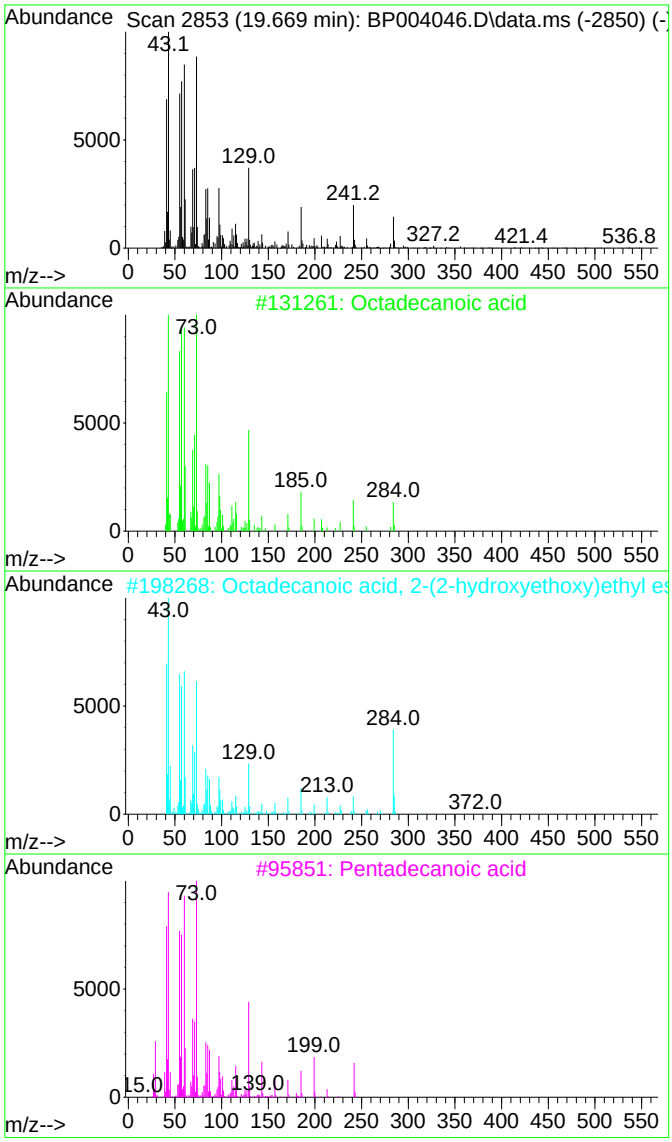
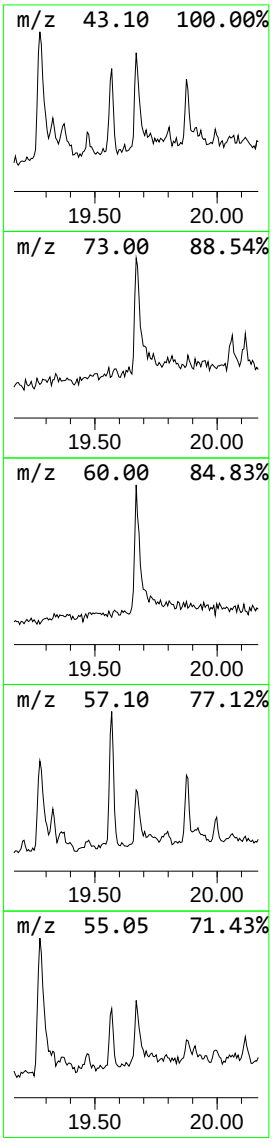
Peak Number 11 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.669	3.00 ng	184744	Chrysene-d12	21.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	74





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Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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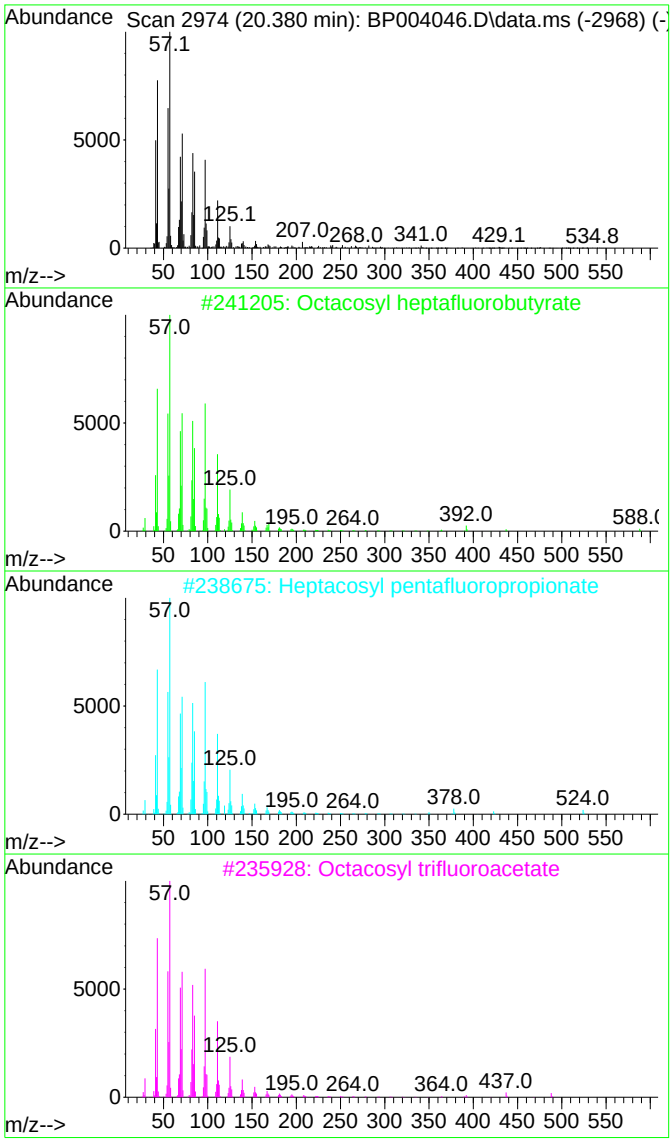
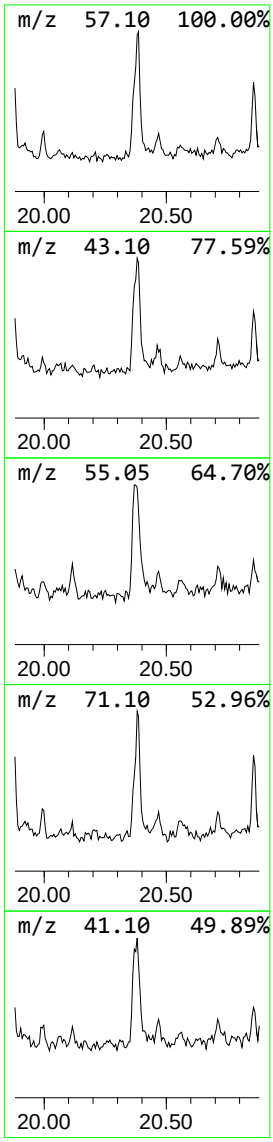
Peak Number 12 Octacosyl heptafluorobutyrate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.380	4.32 ng	266586	Chrysene-d12	21.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosyl heptafluorobutyrate	606	C32H57F7O2	1000351-83-6	91
2	Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91
3	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
4	Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
5	1-Octadecene	252	C18H36	000112-88-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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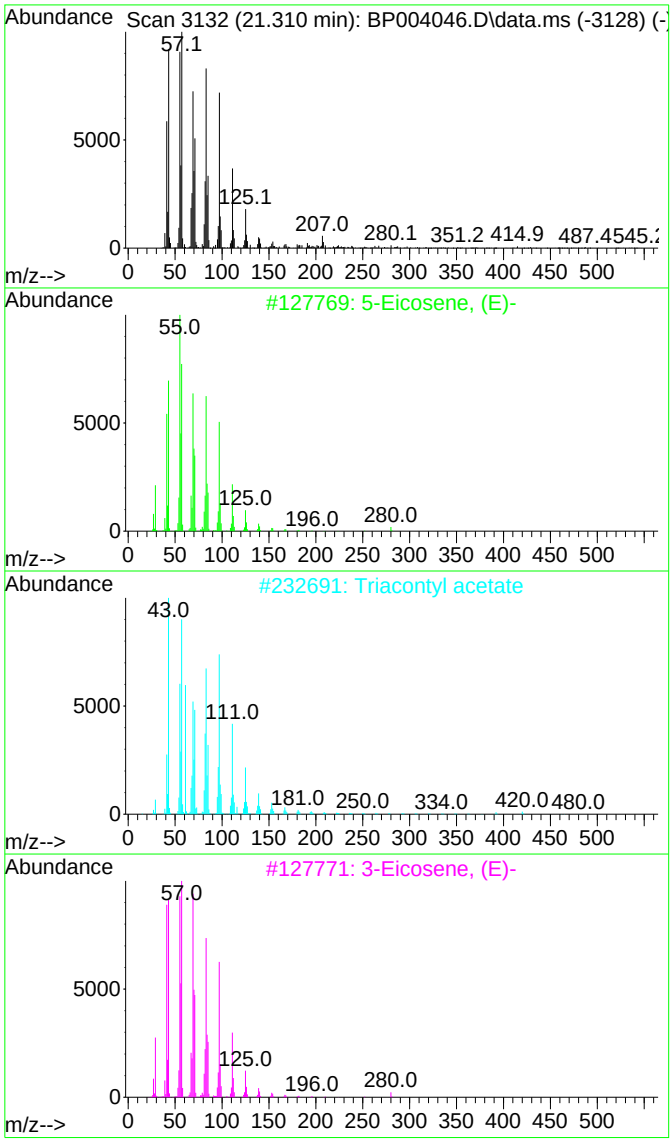
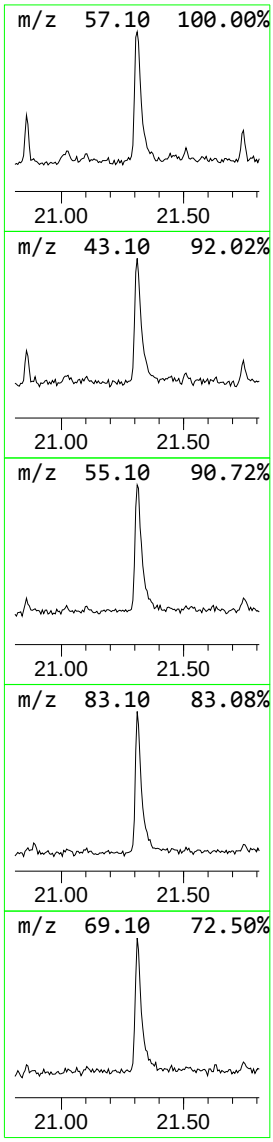
Peak Number 13 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.310	7.61 ng	469525	Chrysene-d12	21.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	93
2	Triacontyl acetate	480	C32H64O2	041755-58-2	91
3	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112120\
Data File : BP004046.D
Acq On : 21 Nov 2020 19:47
Operator : CG/JU
Sample : L4839-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B5-GW

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Cyclohexane	3.075	11.4	ng	364439	1	8.263	641927	20.0
Butane, 2-metho...	3.164	37.0	ng	1187530	1	8.263	641927	20.0
Methyl Isobutyl...	3.840	2.7	ng	85753	1	8.263	641927	20.0
2-Pentanone, 4-...	5.264	4.5	ng	142915	1	8.263	641927	20.0
unknown6.399	6.399	2.4	ng	75969	1	8.263	641927	20.0
2,4,7,9-Tetrame...	13.775	9.6	ng	512806	3	14.887	1063820	20.0
unknown15.986	15.986	6.1	ng	324327	3	14.887	1063820	20.0
n-Hexadecanoic ...	18.463	11.7	ng	717027	4	17.616	1223560	20.0
Cetene	19.275	5.0	ng	303220	4	17.616	1223560	20.0
Diisooctyl maleate	19.569	2.8	ng	173215	4	17.616	1223560	20.0
Octadecanoic acid	19.669	3.0	ng	184744	5	21.651	1233430	20.0
Octacosyl hepta...	20.380	4.3	ng	266586	5	21.651	1233430	20.0
5-Eicosene, (E)-	21.310	7.6	ng	469525	5	21.651	1233430	20.0