LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

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

Signal : TIC: BP003837.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.952	5	11	29	rVV	6001524	9029072	100.00%	14.628%
2	3.105	30	37	47	rVV3	126040	321544	3.56%	0.521%
3	3.199	47	53	67	rVB	389506	627214	6.95%	1.016%
4	5.822	490	499	510	rBV	3075029	4717847	52.25%	7.643%
5	7.469	771	779	793	rBV	3002246	4939977	54.71%	8.003%
6	7.869	841	847	856	rVB	97065	146688	1.62%	0.238%
7	8.305	913	921	929	rBV	603664	972141	10.77%	1.575%
8	9.469	1110	1119	1130	rVB	1829650	3109307	34.44%	5.037%
9	11.134	1395	1402	1410	rBV	767146	1310797	14.52%	2.124%
10	13.546	1805	1812	1819	rBV	4508250	6497496	71.96%	10.526%
11	14.346	1942	1948	1957	rVB	591291	793931	8.79%	1.286%
12	14.646	1988	1999	2001	rBV3	81269	236200	2.62%	0.383%
13	14.698	2005	2008	2022	rVB	116060	230477	2.55%	0.373%
14	14.922	2040	2046	2052	rVB	1172912	1686191	18.68%	2.732%
15	16.398	2290	2297	2311	rBV	3454175	5043430	55.86%	8.171%
16	17.651	2503	2510	2514	rBV	1365613	1978142	21.91%	3.205%
17	17.692	2514	2517	2523	rVV	319223	430768	4.77%	0.698%
18	18.487	2646	2652	2658	rBV2	722627	1024823	11.35%	1.660%
19	18.539	2658	2661	2669	rVB2	133549	206088	2.28%	0.334%
20	18.681	2680	2685	2689	rBV2	84910	169149	1.87%	0.274%
21	18.998	2736	2739	2751	rVB2	57875	107755	1.19%	0.175%
22	19.369	2797	2802	2806	rBV	87796	131088	1.45%	0.212%
23	19.622	2840	2845	2850	rVB	389014	522959	5.79%	0.847%
24	19.686	2853	2856	2864	rVB2	170439	233669	2.59%	0.379%
25	19.969	2900	2904	2911	rVB	428024	576486	6.38%	0.934%
26	20.145	2928	2934	2939	rBV	6700148	8334294	92.31%	13.502%
27	20.610	3008	3013	3019	rVB2	94426	202688	2.24%	0.328%
28	20.769	3037	3040	3044	rVB2	108817	126521	1.40%	0.205%
29	21.322	3131	3134	3143	rVB2	946700	1262915	13.99%	2.046%
30	21.398	3143	3147	3151	rBV	319363	366582	4.06%	0.594%
31	21.675	3187	3194	3198	rBV	1478680	2218898	24.58%	3.595%
32	21.710	3198	3200	3205	rVB	193045	211108	2.34%	0.342%
33	22.257	3287	3293	3300	rVB3	184634	378967	4.20%	0.614%
34	22.898	3399	3402	3412	rVB	119304	186897	2.07%	0.303%
35	22.992	3415	3418	3426	rVB3	170954	257575	2.85%	0.417%
36	23.304	3467	3471	3475	rBV	138970	192174	2.13%	0.311%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

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

37	23.351	3476	3479	3483	rBV2	149171	229124	2.54%	0.371%
38	23.933	3575	3578	3587	rVB	188108	338421	3.75%	0.548%
39	24.151	3609	3615	3622	rBV	923802	1906921	21.12%	3.089%
40	24.745	3712	3716	3730	rVB3	194815	469434	5.20%	0.761%

Sum of corrected areas: 61725758

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
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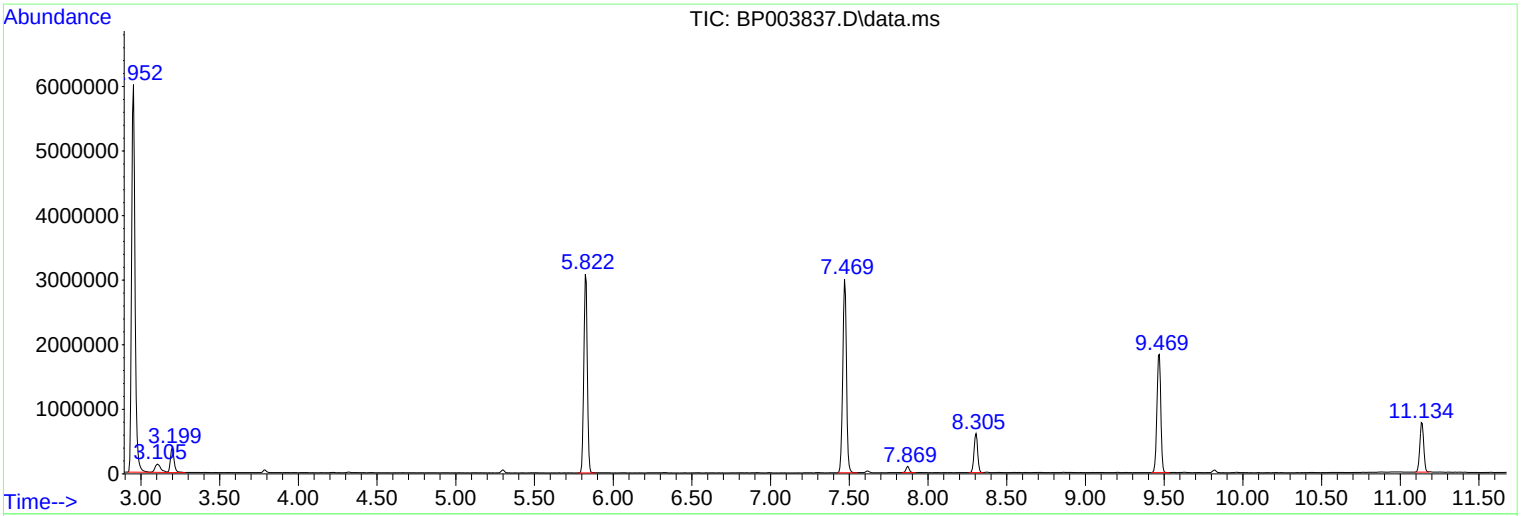
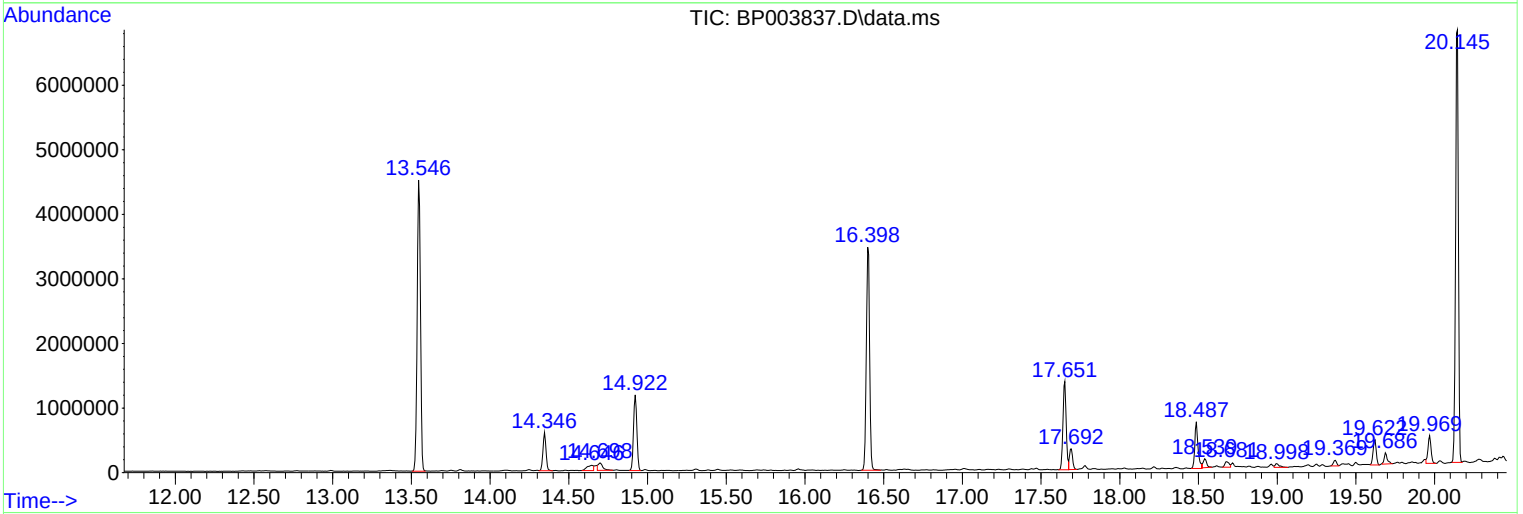
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

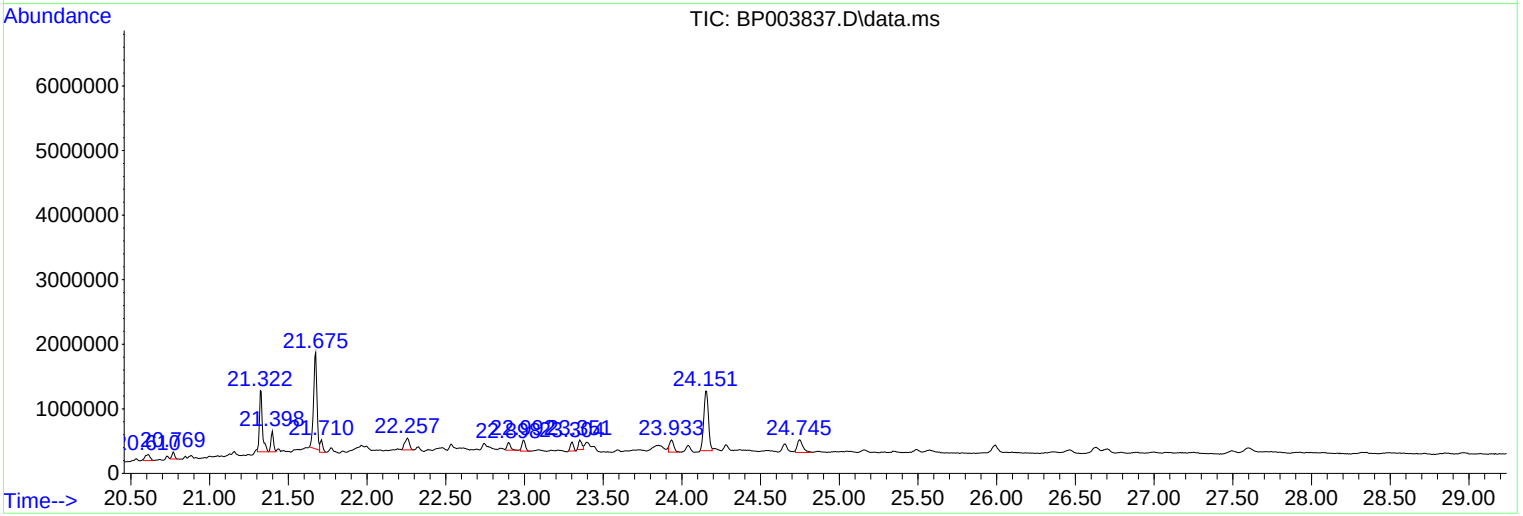
Instrument :
BNA_P
ClientSampleId :
356

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

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

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

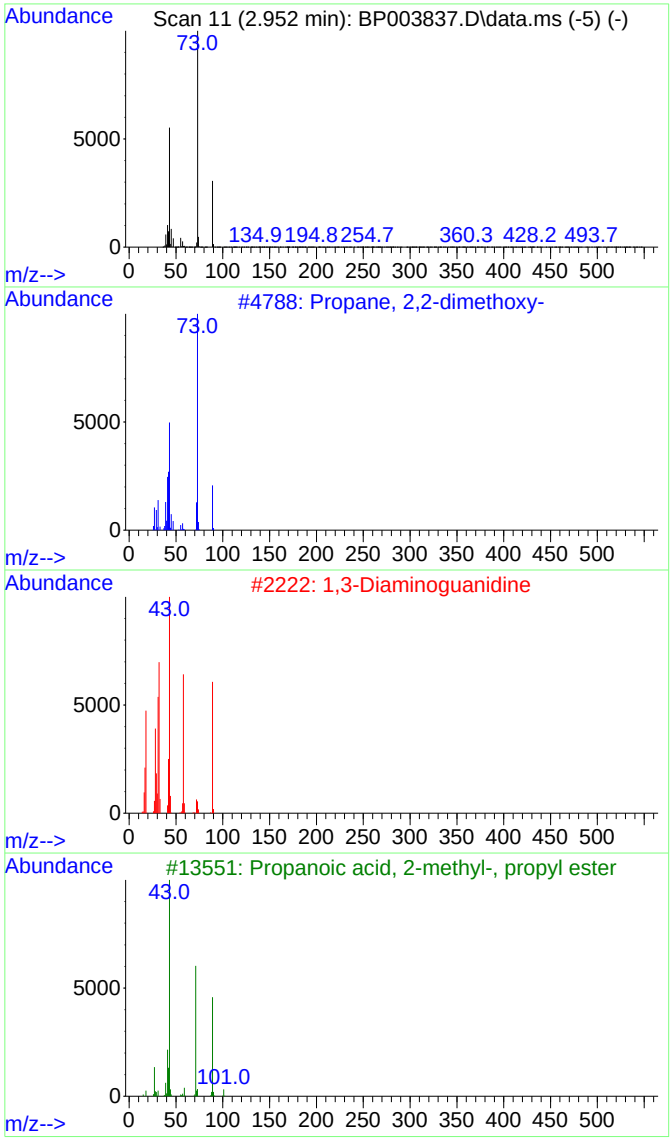
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

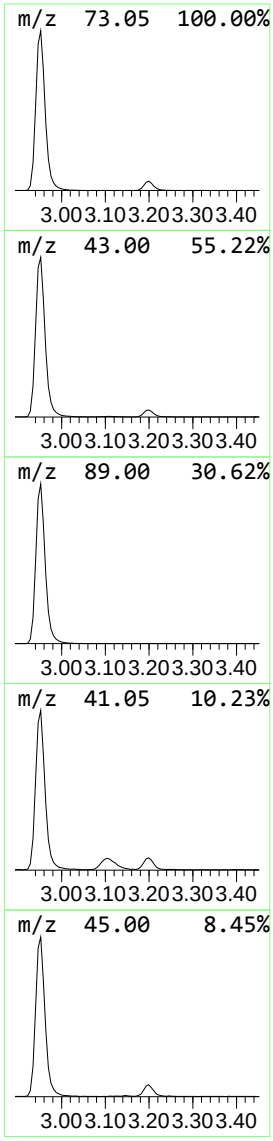
****

Peak Number 1 unknown2.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.952	185.76 ng	9029070	1,4-Dichlorobenzene-d4	8.305

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2	1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
3	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
4	1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

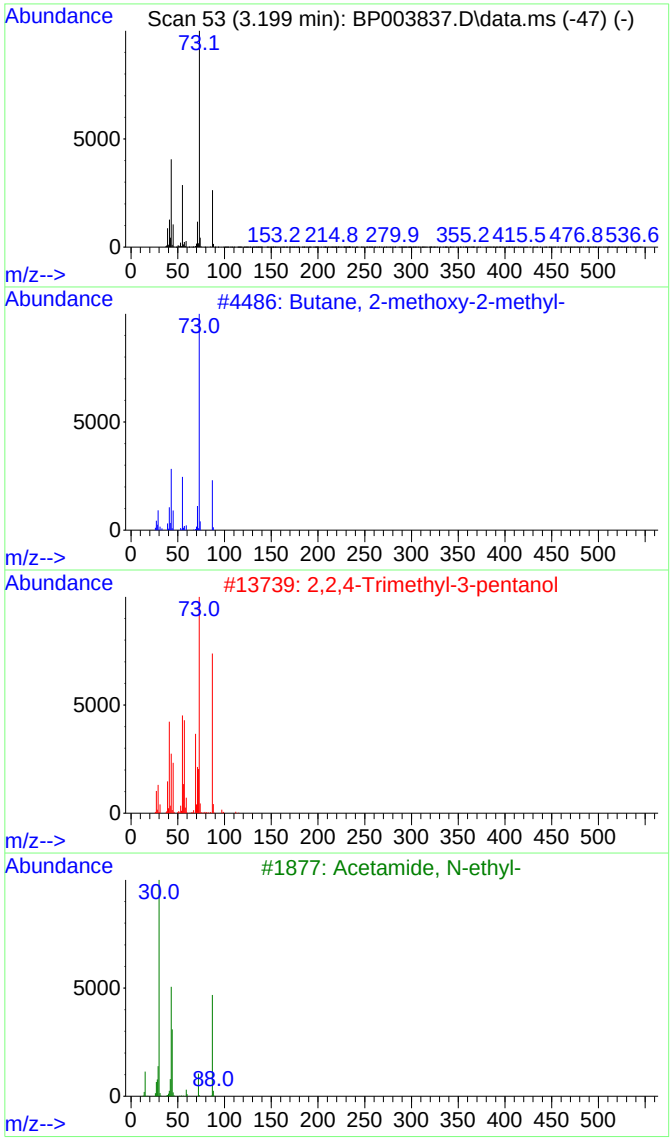
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

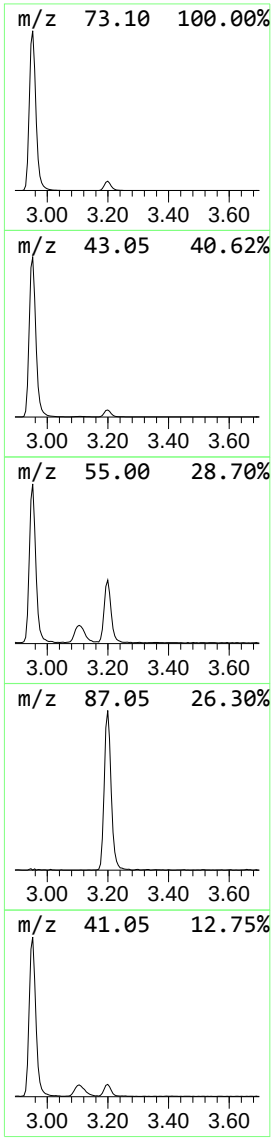
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.199	12.90 ng	627214	1,4-Dichlorobenzene-d4	8.305

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	39
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	17
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

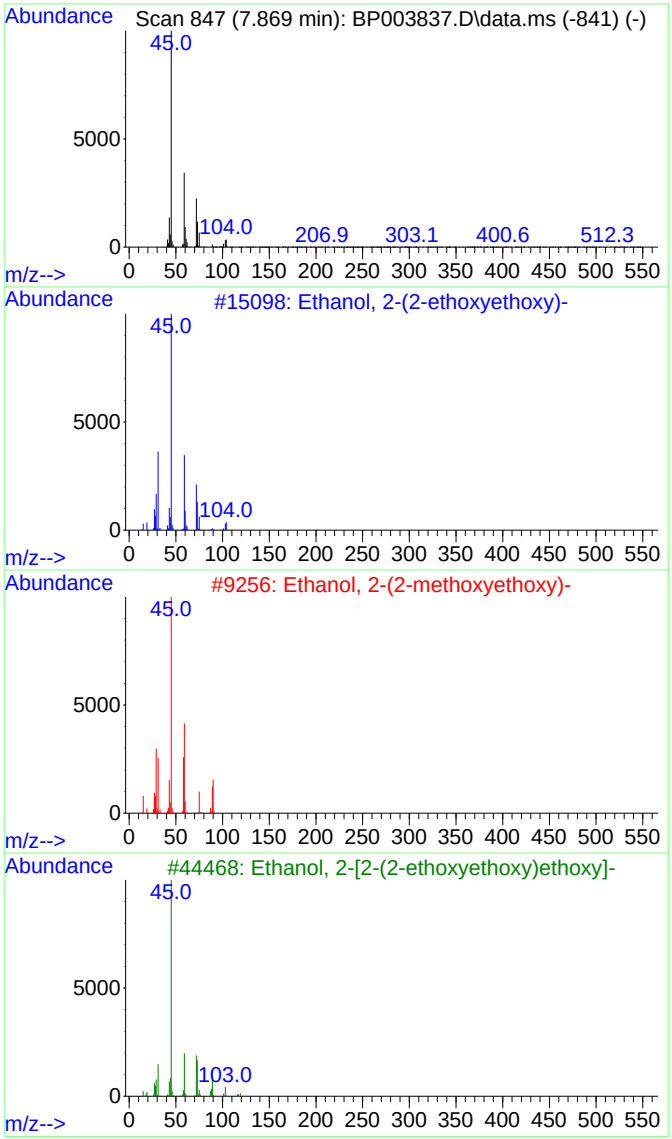
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

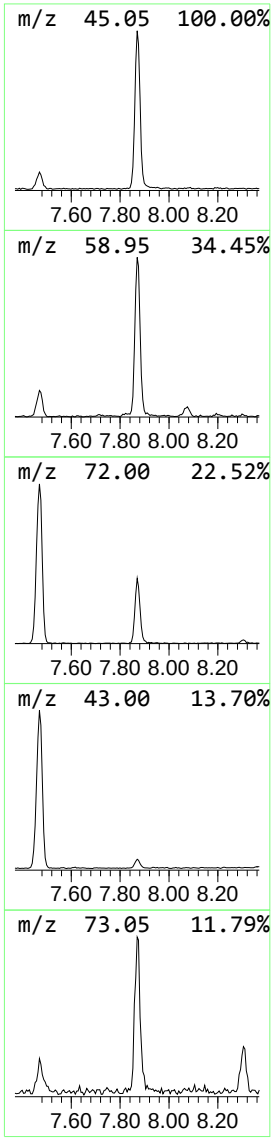
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.869	3.02 ng	146688	1,4-Dichlorobenzene-d4	8.305

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2	Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	42
3	Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	39
4	Diethyl carbitol	162	C8H18O3	000112-36-7	33
5	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

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

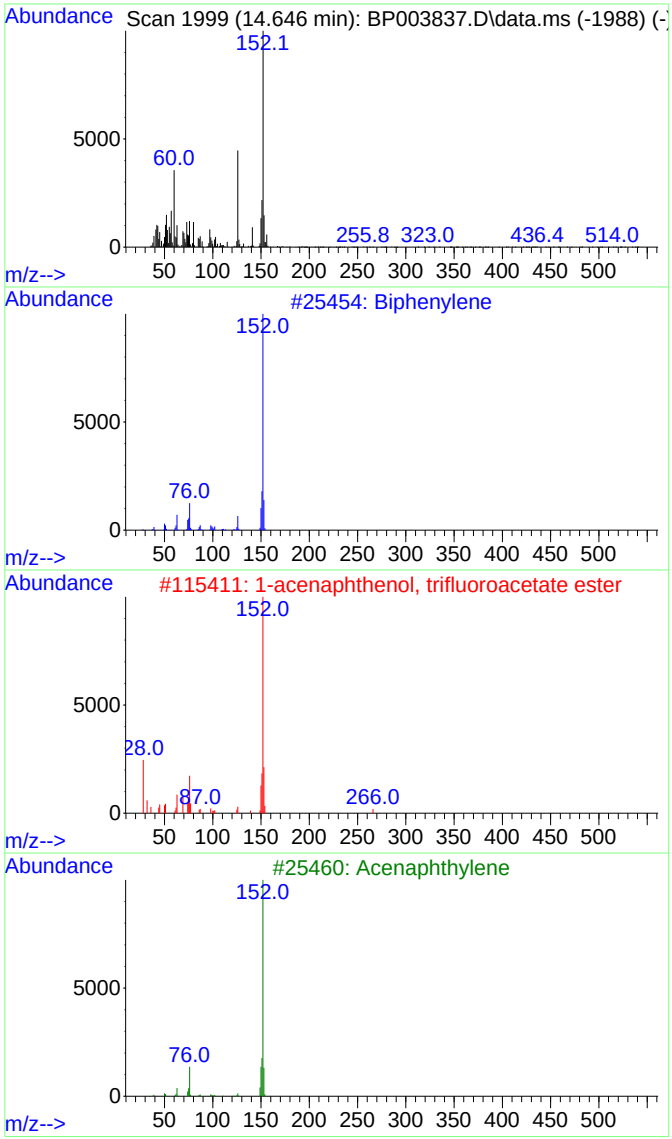
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

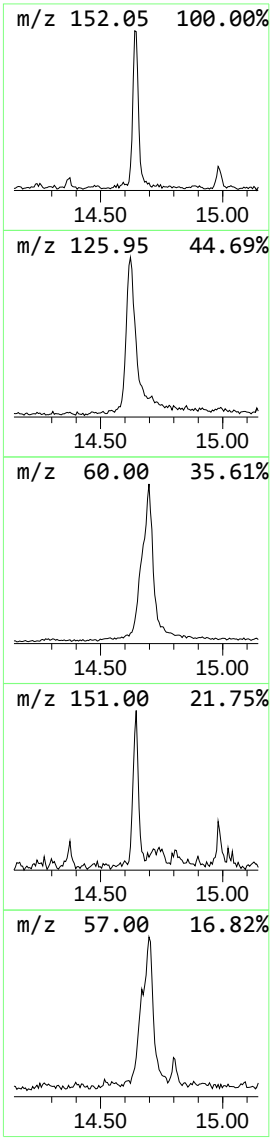
****

Peak Number 5 Biphenylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.646	2.80 ng	236200	Acenaphthene-d10	14.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenylene	152	C12H8	000259-79-0	60
2	1-acenaphthenol, trifluoroacetat...	266	C14H9F3O2	1000376-45-2	43
3	Acenaphthylene	152	C12H8	000208-96-8	43
4	Benzaldehyde, 2-hydroxy-5-methoxy-	152	C8H8O3	000672-13-9	38
5	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

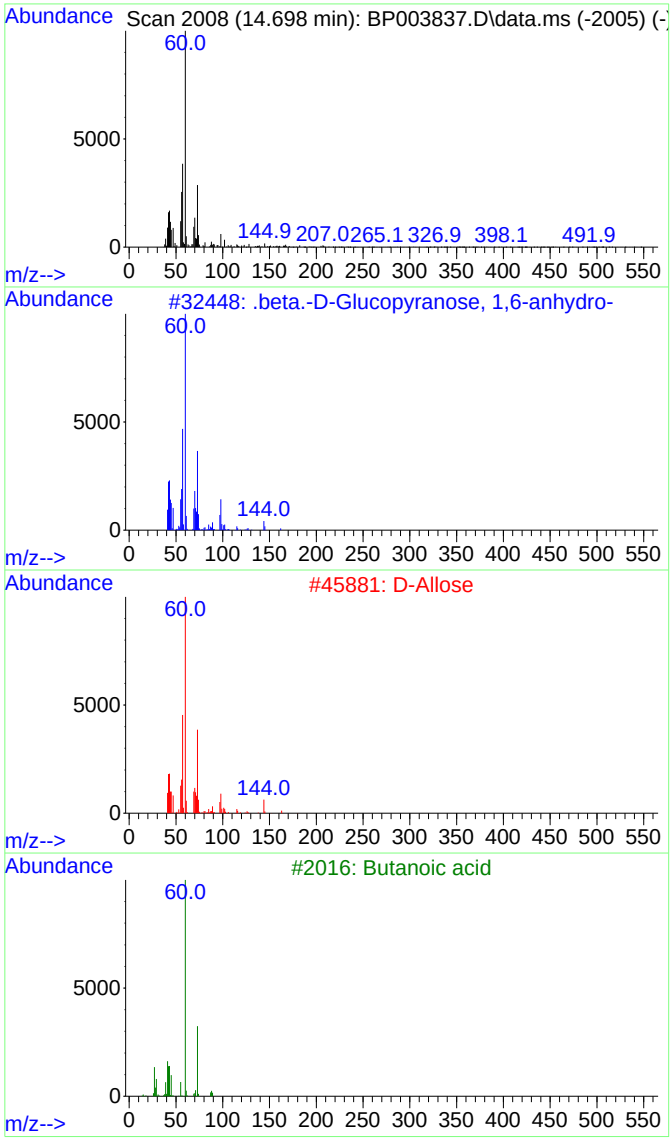
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

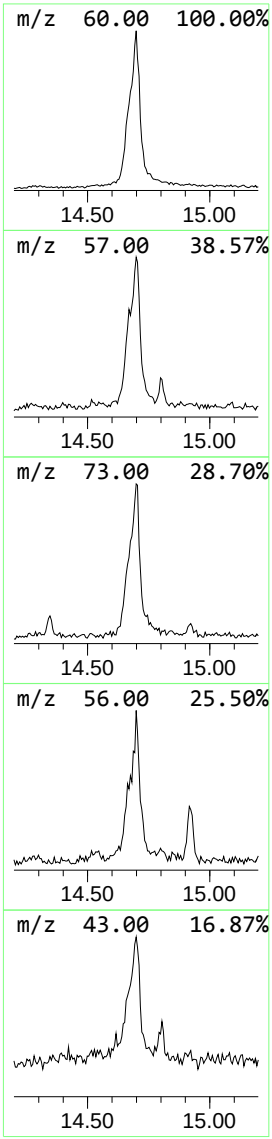
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 .beta.-D-Glucopyranose, 1,6... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.698	2.73 ng	230477	Acenaphthene-d10	14.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-D-Glucopyranose, 1,6-anhy...	162	C6H10O5	000498-07-7	59
2	D-Allose	180	C6H12O6	002595-97-3	56
3	Butanoic acid	88	C4H8O2	000107-92-6	50
4	Pentanoic acid	102	C5H10O2	000109-52-4	40
5	Heptanoic acid	130	C7H14O2	000111-14-8	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

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

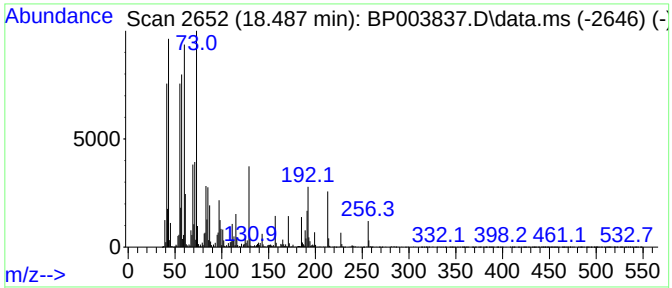
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

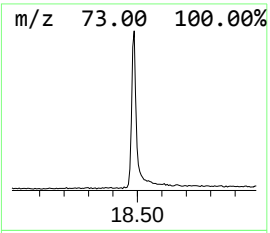
****

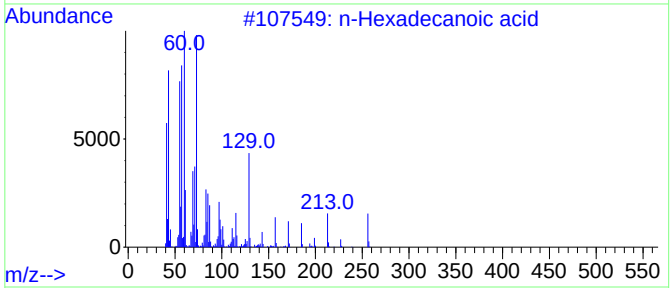
Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

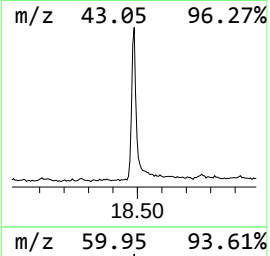
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	10.36 ng	1024820	Phenanthrene-d10	17.651

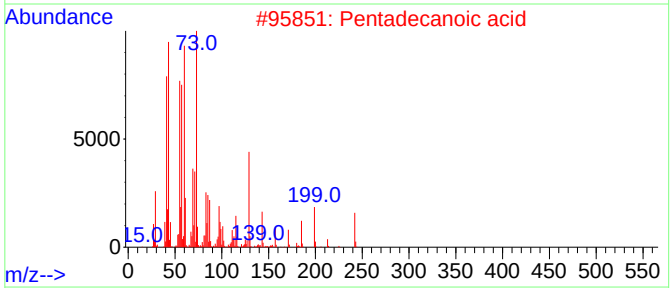
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Octadecanoic acid	284	C18H36O2	000057-11-4	91
4	Tridecanoic acid	214	C13H26O2	000638-53-9	90
5	Undecanoic acid	186	C11H22O2	000112-37-8	76

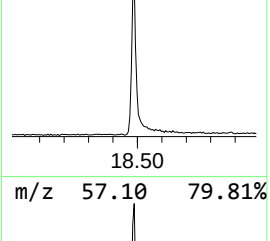


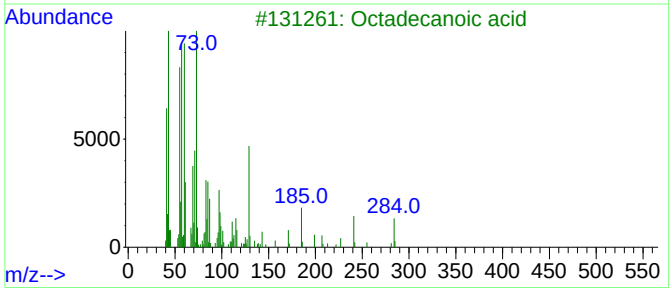


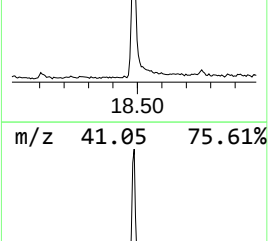


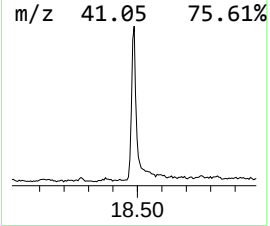












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

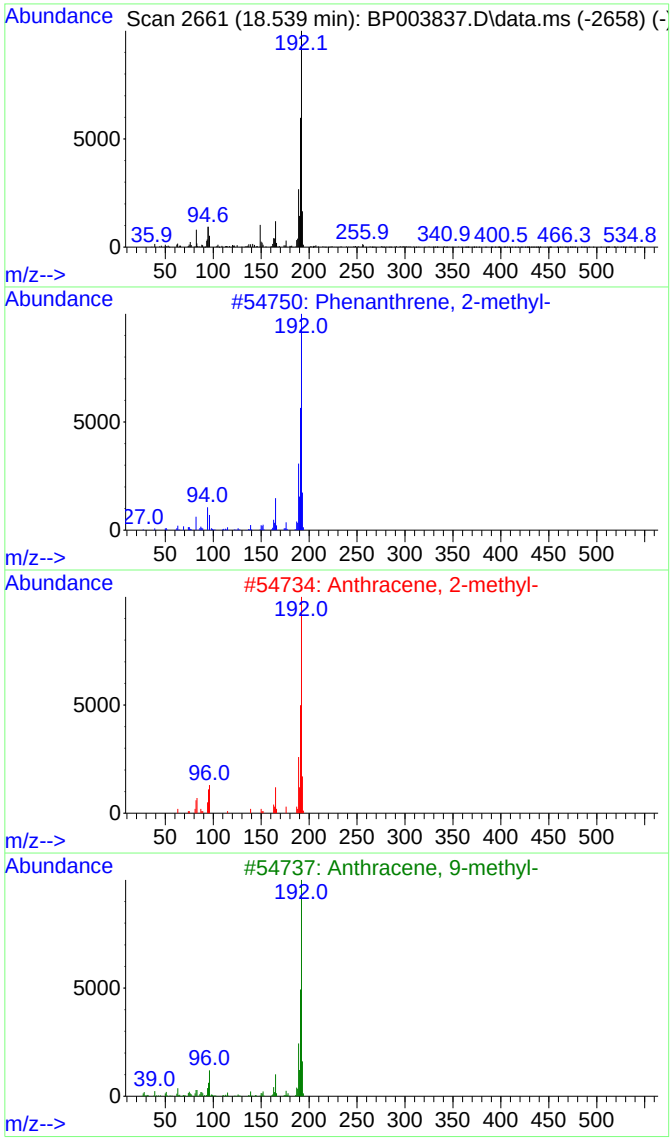
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

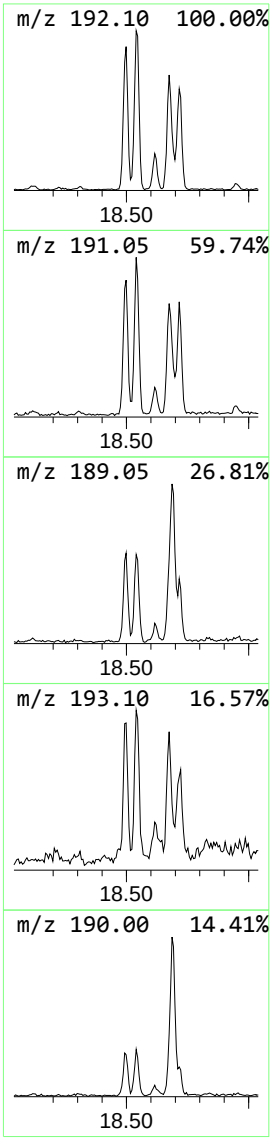
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.540	2.08 ng	206088	Phenanthrene-d10	17.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89





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Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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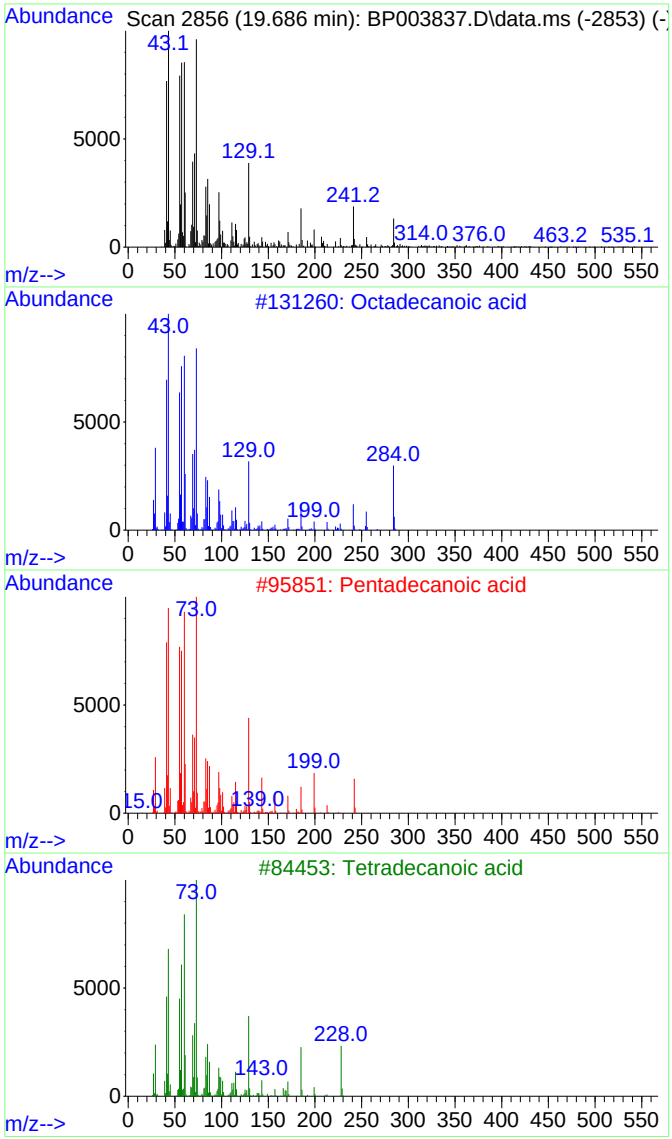
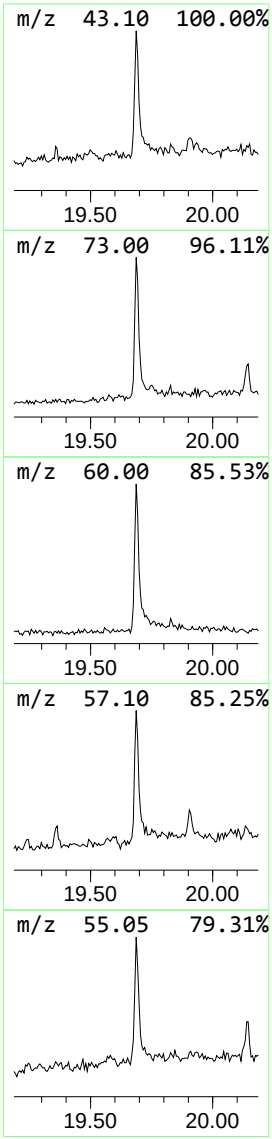
Peak Number 9 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.686	2.11 ng	233669	Chrysene-d12	21.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	83
5	n-Decanoic acid	172	C10H20O2	000334-48-5	64





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Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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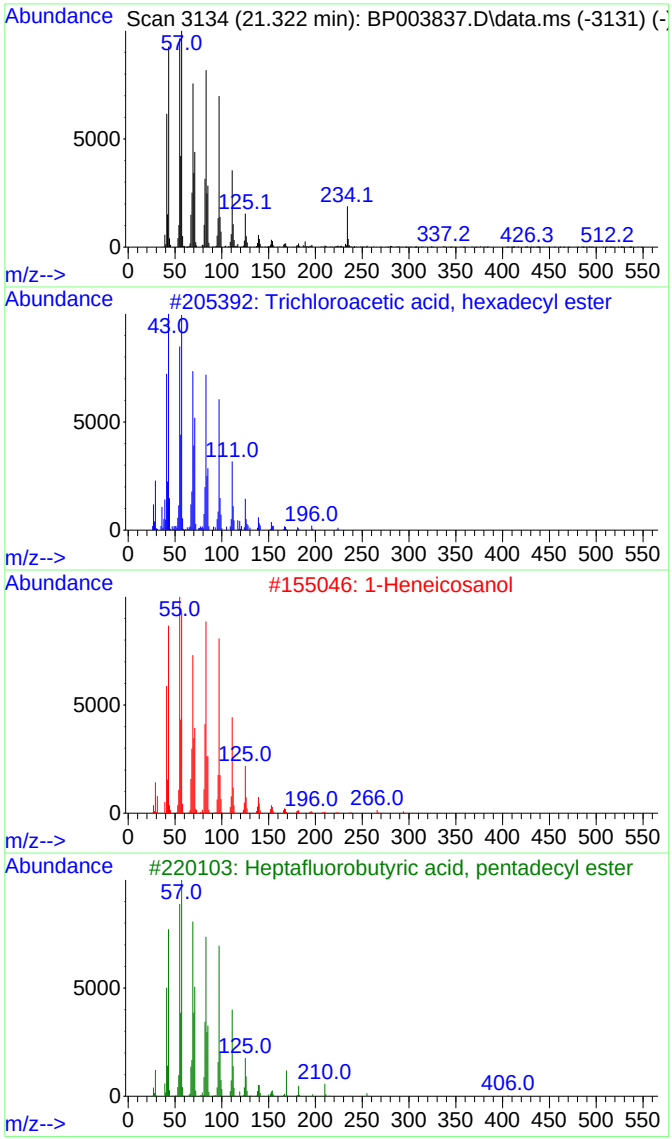
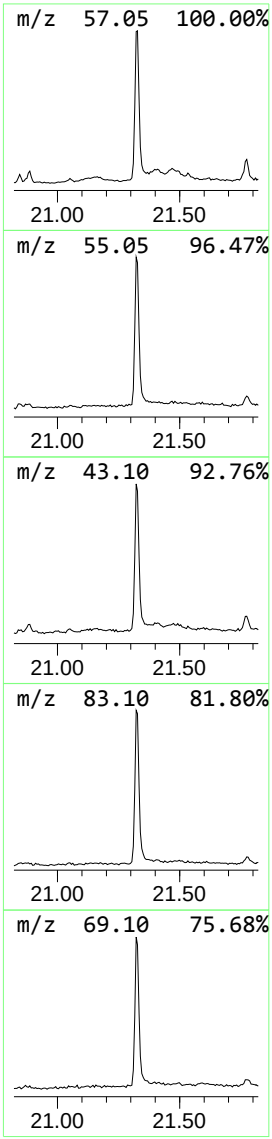
Peak Number 10 Trichloroacetic acid, hexad... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.322	11.38 ng	1262920	Chrysene-d12	21.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91





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Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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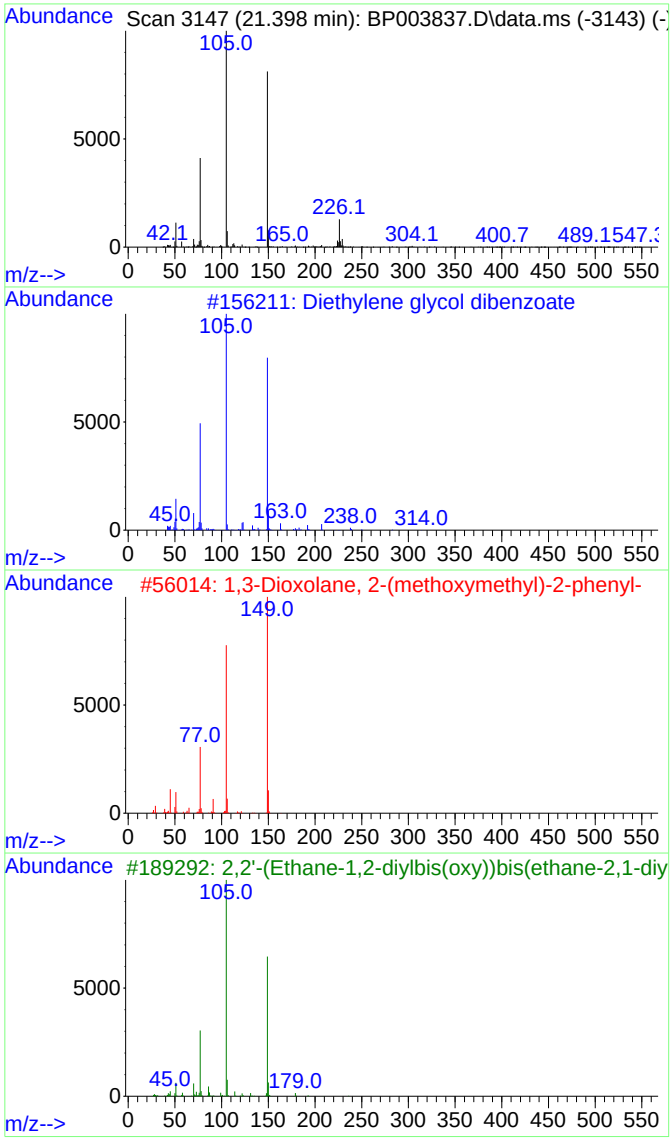
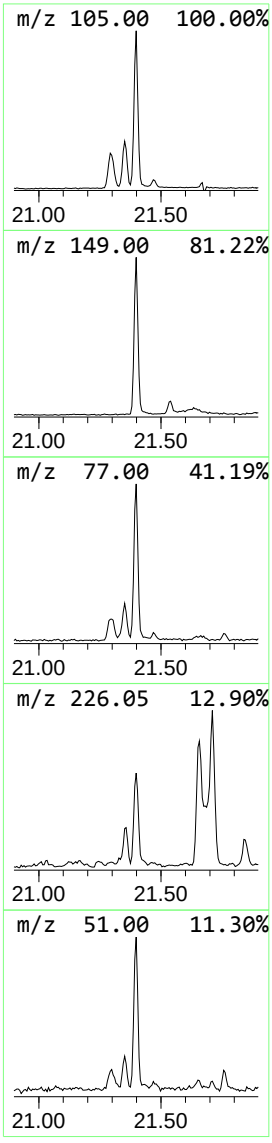
Peak Number 11 Diethylene glycol dibenzoate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.398	3.30 ng	366582	Chrysene-d12	21.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	80
2	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	72
3	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	56
4	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	56
5	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	50





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Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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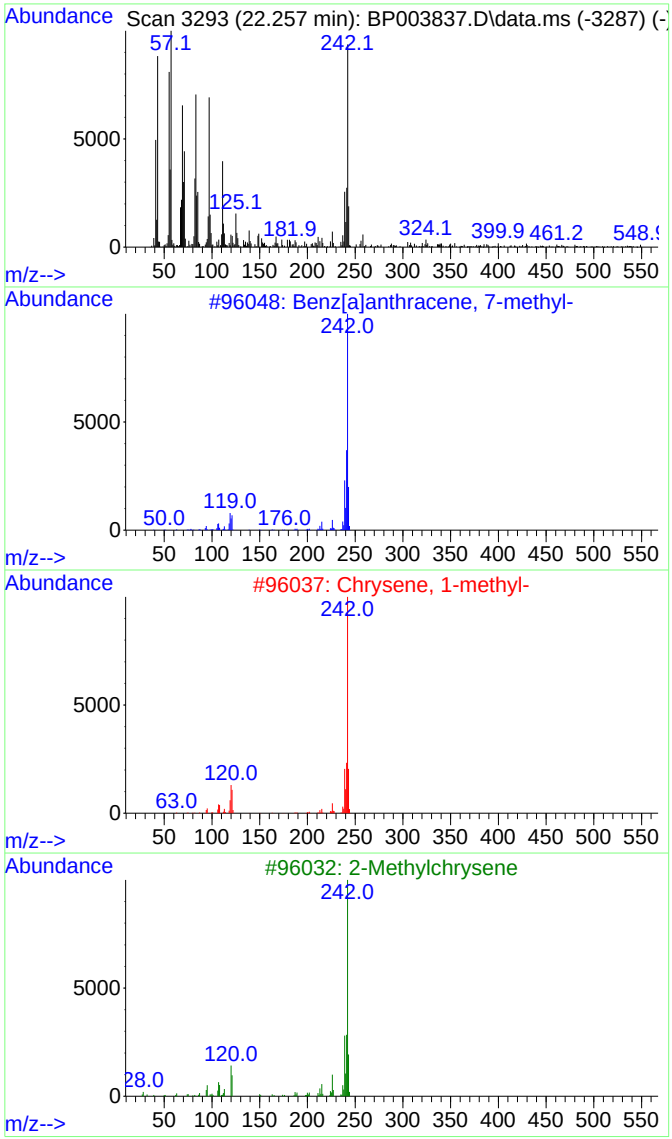
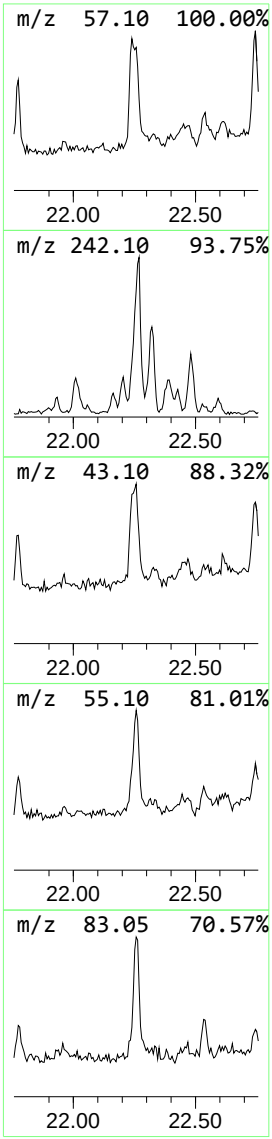
Peak Number 12 Benz[a]anthracene, 7-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.257	3.42 ng	378967	Chrysene-d12	21.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	52
2	Chrysene, 1-methyl-	242	C19H14	003351-28-8	50
3	2-Methylchrysene	242	C19H14	003351-32-4	47
4	Chrysene, 5-methyl-	242	C19H14	003697-24-3	38
5	Chrysene, 6-methyl-	242	C19H14	001705-85-7	38





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Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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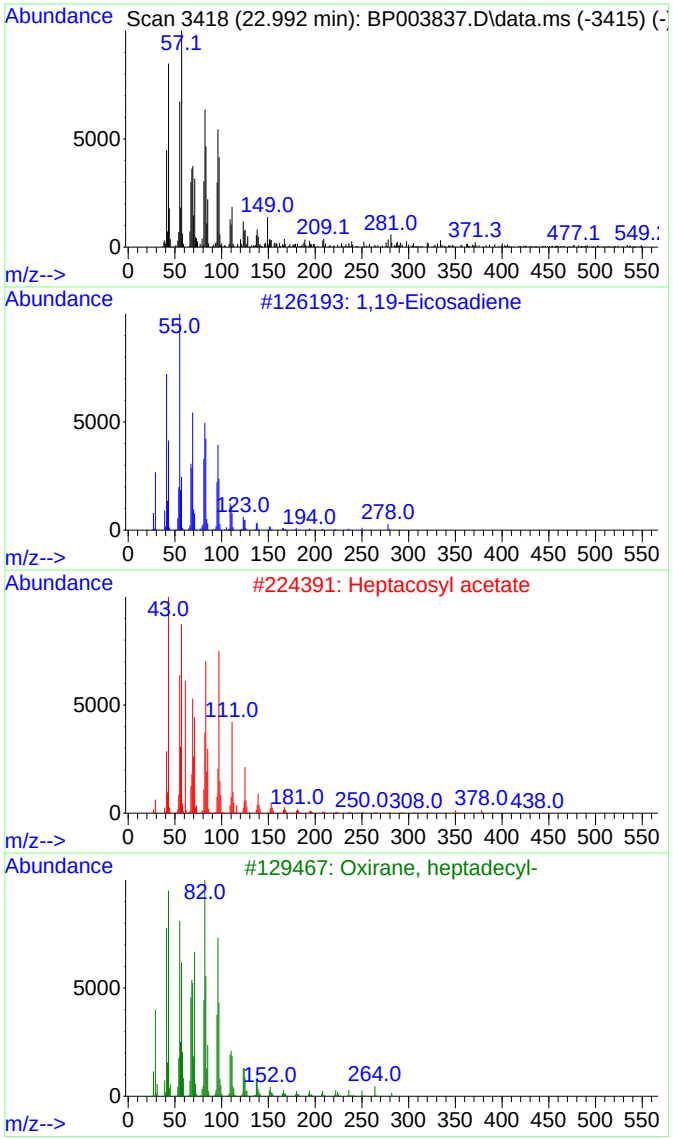
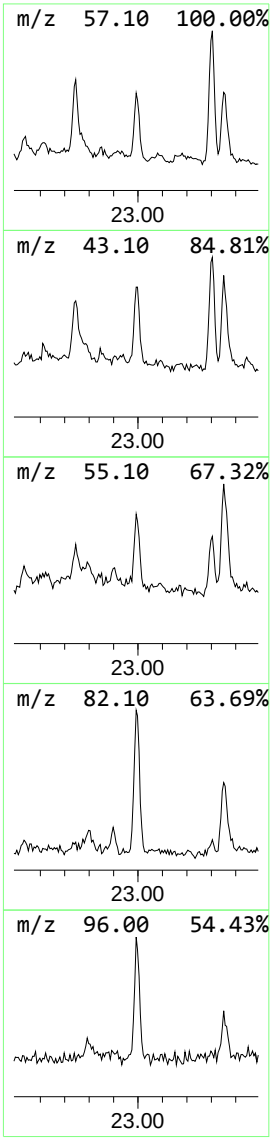
Peak Number 13 1,19-Eicosadiene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.992	2.70 ng	257575	Perylene-d12	24.157

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	91
2	Heptacosyl acetate	438	C29H58O2	1000351-78-2	87
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	68
4	2-Octadecyl-propane-1,3-diol	328	C21H44O2	005337-61-1	64
5	Hexadecanal	240	C16H32O	000629-80-1	62





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Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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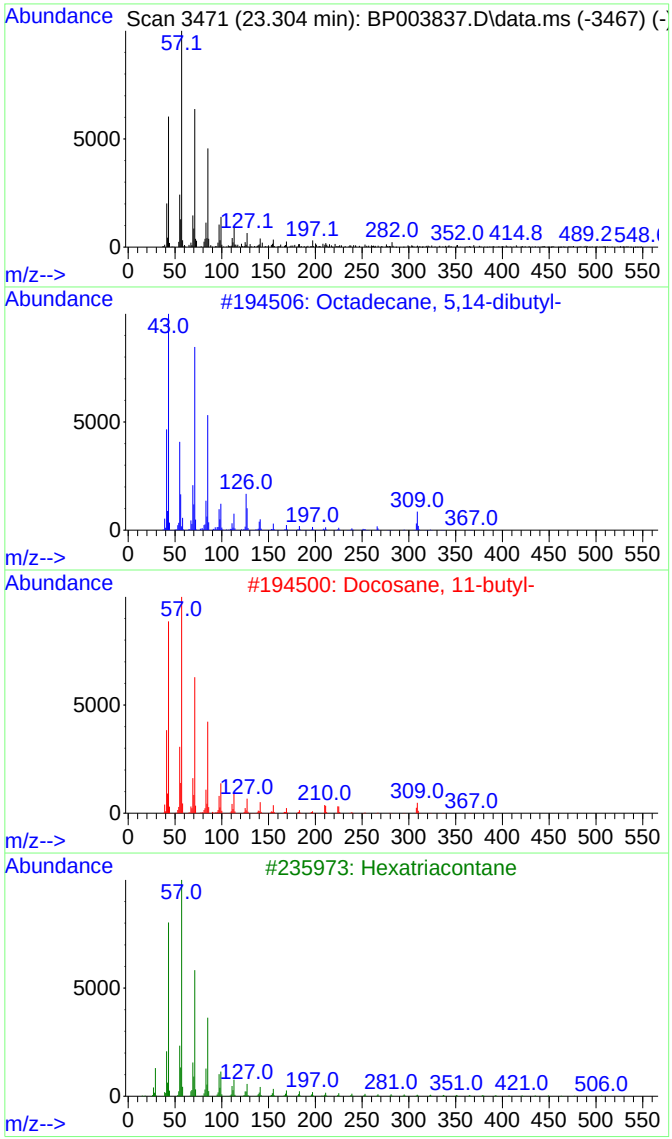
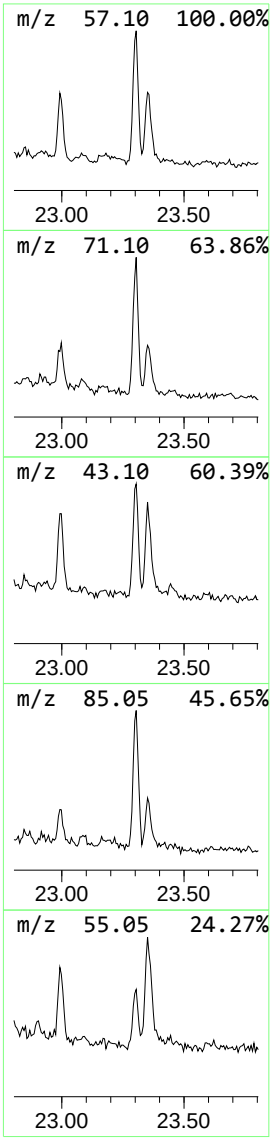
Peak Number 14 Octadecane, 5,14-dibutyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.304	2.02 ng	192174	Perylene-d12	24.157

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	94
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3	Hexatriacontane	507	C36H74	000630-06-8	93
4	Eicosane	282	C20H42	000112-95-8	93
5	Octacosane	394	C28H58	000630-02-4	81





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Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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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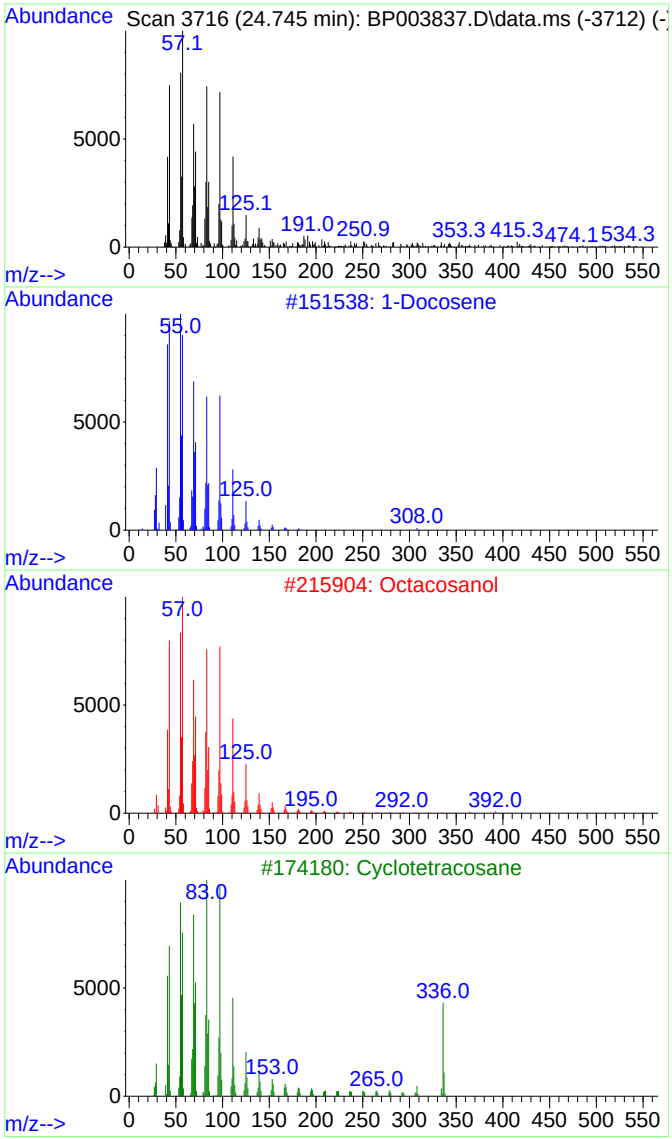
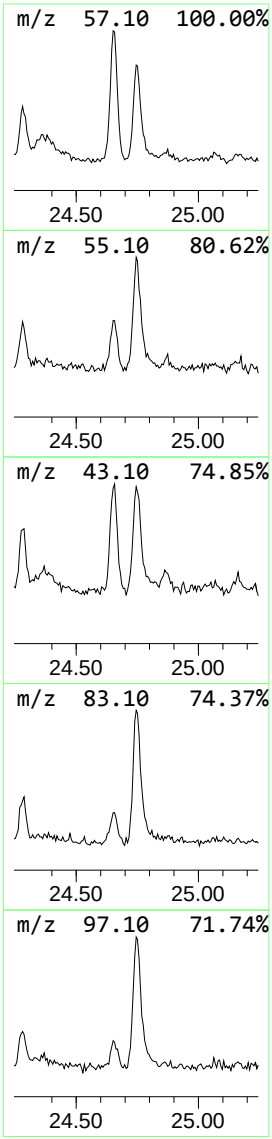
Peak Number 16 1-Docosene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.745	4.92 ng	469434	Perylene-d12	24.157

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	Octacosanol	410	C28H58O	000557-61-9	94
3	Cyclotetracosane	336	C24H48	000297-03-0	91
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5	Heneicosyl trifluoroacetate	408	C23H43F3O2	1000351-76-5	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110620\
Data File : BP003837.D
Acq On : 06 Nov 2020 13:26
Operator : CG/JU
Sample : L4635-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 356

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-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
unknown2.95	2.952	185.8	ng	9029070	1	8.305	972141	20.0
Butane, 2-metho...	3.199	12.9	ng	627214	1	8.305	972141	20.0
Ethanol, 2-(2-e...	7.869	3.0	ng	146688	1	8.305	972141	20.0
Biphenylene	14.646	2.8	ng	236200	3	14.922	1686190	20.0
.beta.-D-Glucop...	14.698	2.7	ng	230477	3	14.922	1686190	20.0
n-Hexadecanoic ...	18.486	10.4	ng	1024820	4	17.651	1978140	20.0
Phenanthrene, 2...	18.540	2.1	ng	206088	4	17.651	1978140	20.0
Octadecanoic acid	19.686	2.1	ng	233669	5	21.675	2218900	20.0
Trichloroacetic...	21.322	11.4	ng	1262920	5	21.675	2218900	20.0
Diethylene glyc...	21.398	3.3	ng	366582	5	21.675	2218900	20.0
Benz[a]anthrace...	22.257	3.4	ng	378967	5	21.675	2218900	20.0
1,19-Eicosadiene	22.992	2.7	ng	257575	6	24.157	1906920	20.0
Octadecane, 5,1...	23.304	2.0	ng	192174	6	24.157	1906920	20.0
1-Docosene	24.745	4.9	ng	469434	6	24.157	1906920	20.0