LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

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-BP011422.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008947.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	4	9	36	rVB	1004261	2099821	3.14%	0.947%
2	5.422	407	414	426	rBV	1486202	2380326	3.56%	1.074%
3	7.010	676	684	702	rBV	943496	1685966	2.52%	0.761%
4	7.828	814	823	829	rBV	797532	1307706	1.96%	0.590%
5	8.992	1013	1021	1033	rBV	2165224	3757690	5.62%	1.695%
6	10.628	1290	1299	1316	rBV	966819	1738144	2.60%	0.784%
7	13.092	1711	1718	1733	rBV	6202079	9569451	14.32%	4.318%
8	14.469	1946	1952	1967	rBV	1521886	2386974	3.57%	1.077%
9	15.968	2201	2207	2223	rVB	5193073	7439618	11.13%	3.357%
10	16.639	2314	2321	2340	rVB4	261058	933616	1.40%	0.421%
11	17.227	2415	2421	2427	rBV	2009313	2866725	4.29%	1.293%
12	18.121	2552	2573	2582	rBV	7641263	32691613	48.93%	14.751%
13	18.257	2582	2596	2607	rVV2	7866135	32842800	49.15%	14.819%
14	18.410	2607	2622	2638	rVV	19658346	66818592	100.00%	30.149%
15	18.562	2639	2648	2660	rVB	2114952	5491665	8.22%	2.478%
16	19.245	2761	2764	2771	rVV	615640	1289617	1.93%	0.582%
17	19.368	2773	2785	2819	rVB	7681587	14953449	22.38%	6.747%
18	19.792	2851	2857	2862	rBV	10810871	12589116	18.84%	5.680%
19	20.186	2918	2924	2935	rVB	564342	1313513	1.97%	0.593%
20	20.909	3042	3047	3054	rVB	1178015	1730562	2.59%	0.781%
21	21.321	3112	3117	3125	rBV	2700424	3507776	5.25%	1.583%
22	21.556	3152	3157	3165	rBV	1402271	2324257	3.48%	1.049%
23	22.268	3273	3278	3288	rVB	1337824	2321198	3.47%	1.047%
24	23.056	3406	3412	3421	rVB	1135516	2315956	3.47%	1.045%
25	23.727	3520	3526	3535	rVB2	1778872	3604266	5.39%	1.626%
26	23.962	3560	3566	3577	rVB	729034	1668974	2.50%	0.753%

Sum of corrected areas: 221629391

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
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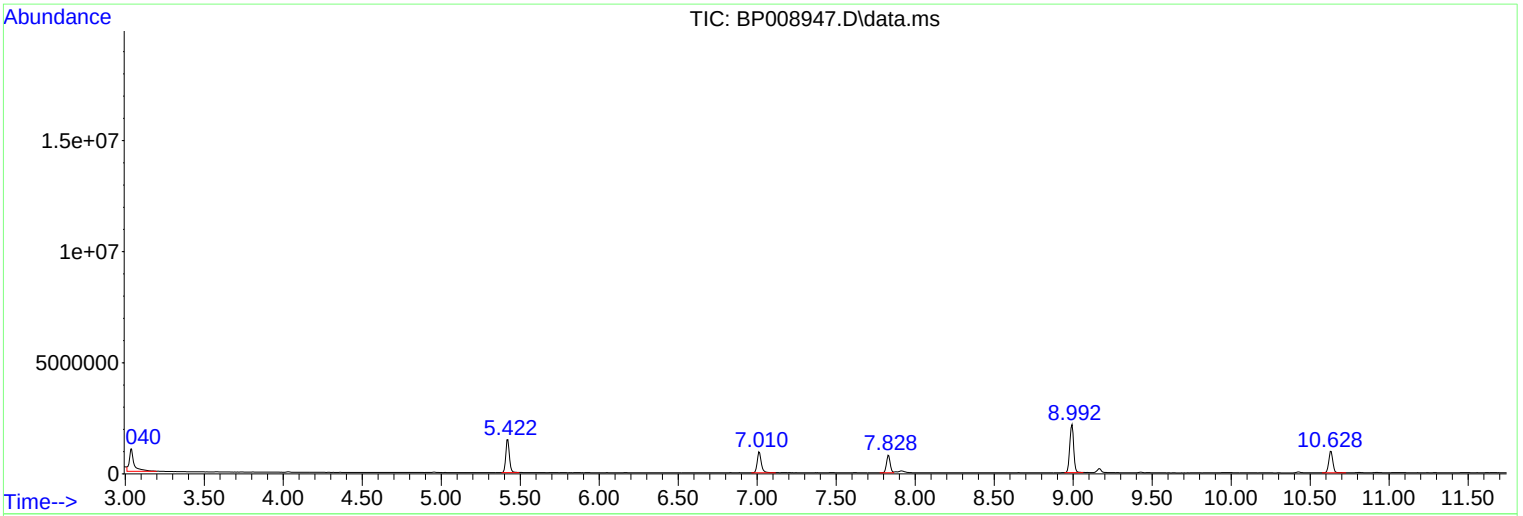
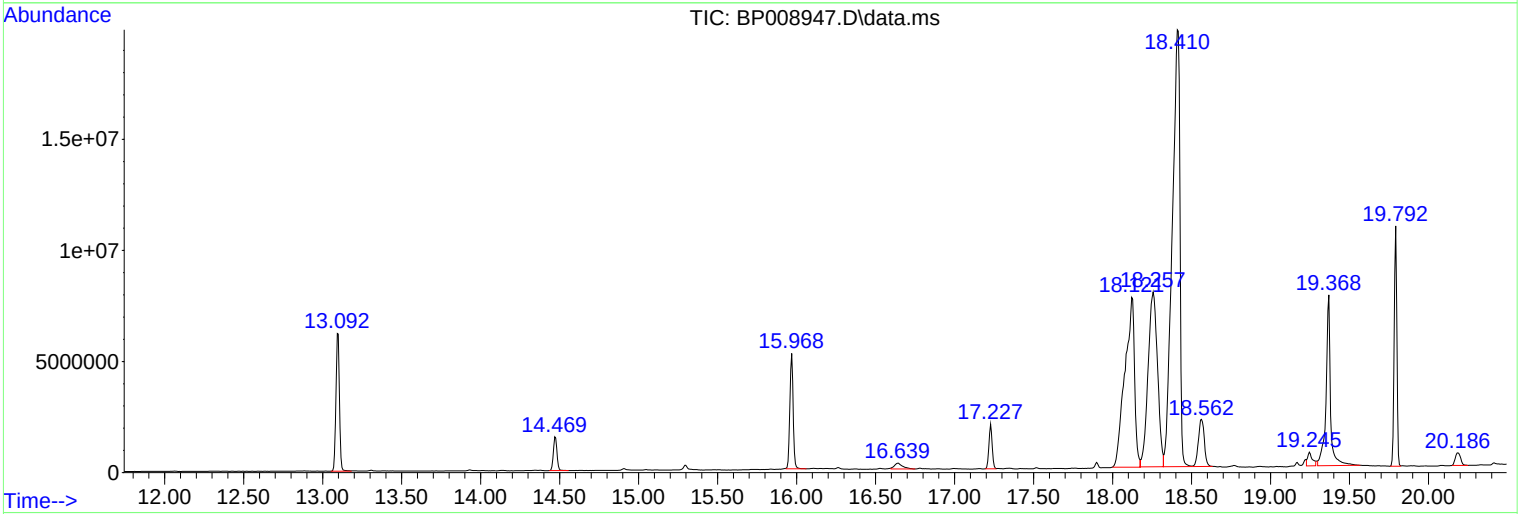
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

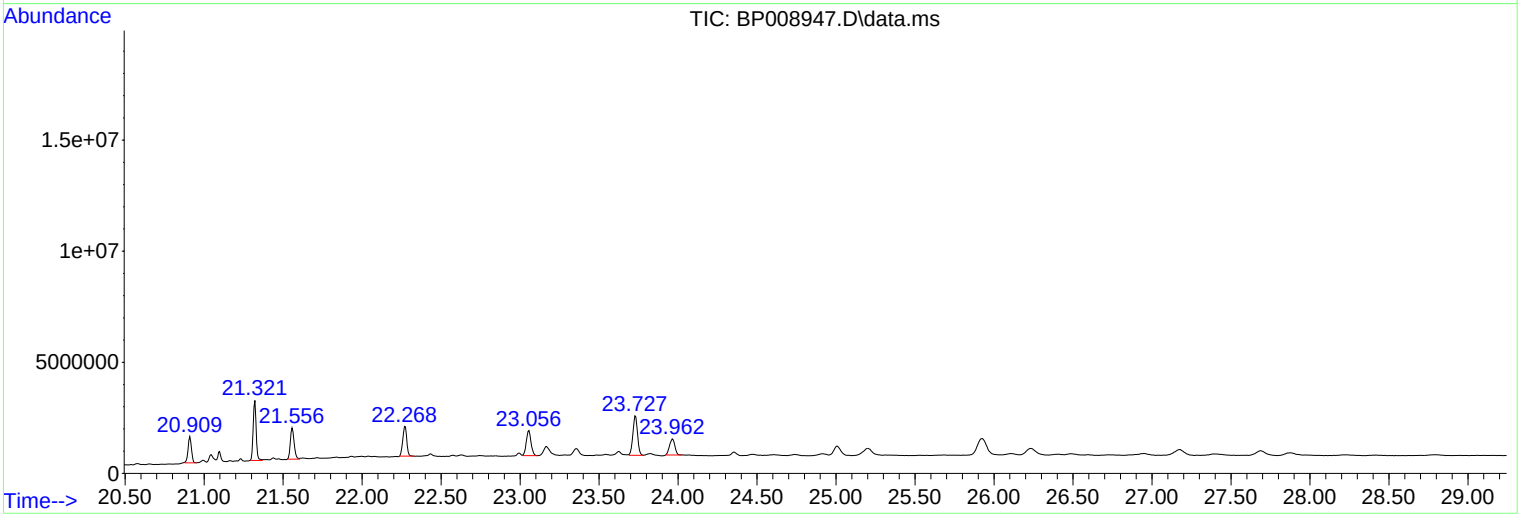
Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

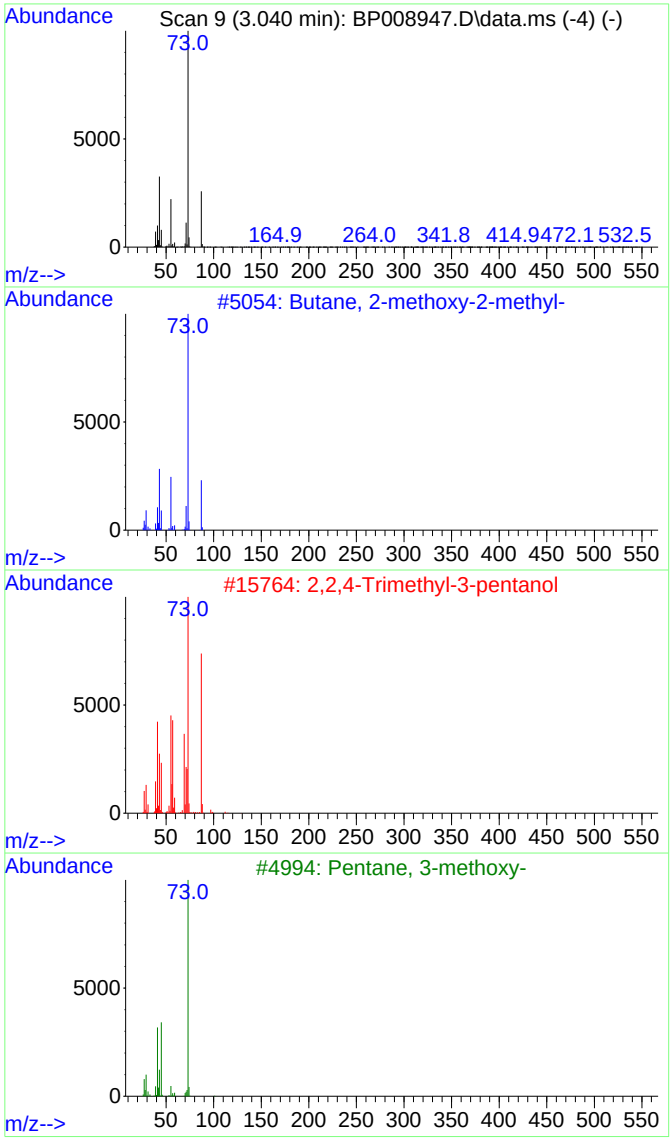
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

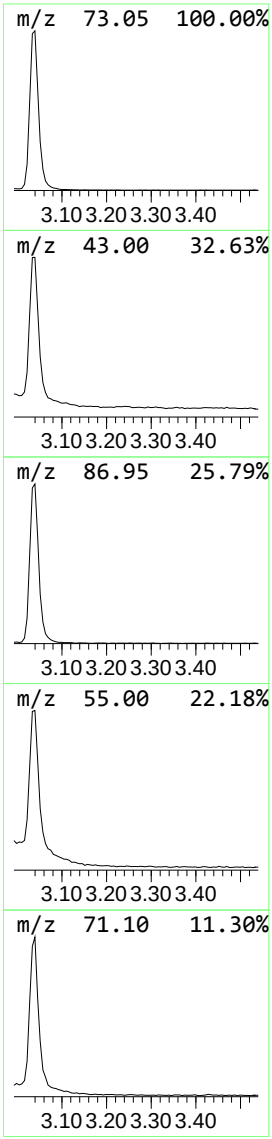
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	32.11 ng	2099820	1,4-Dichlorobenzene-d4	7.828

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	43
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

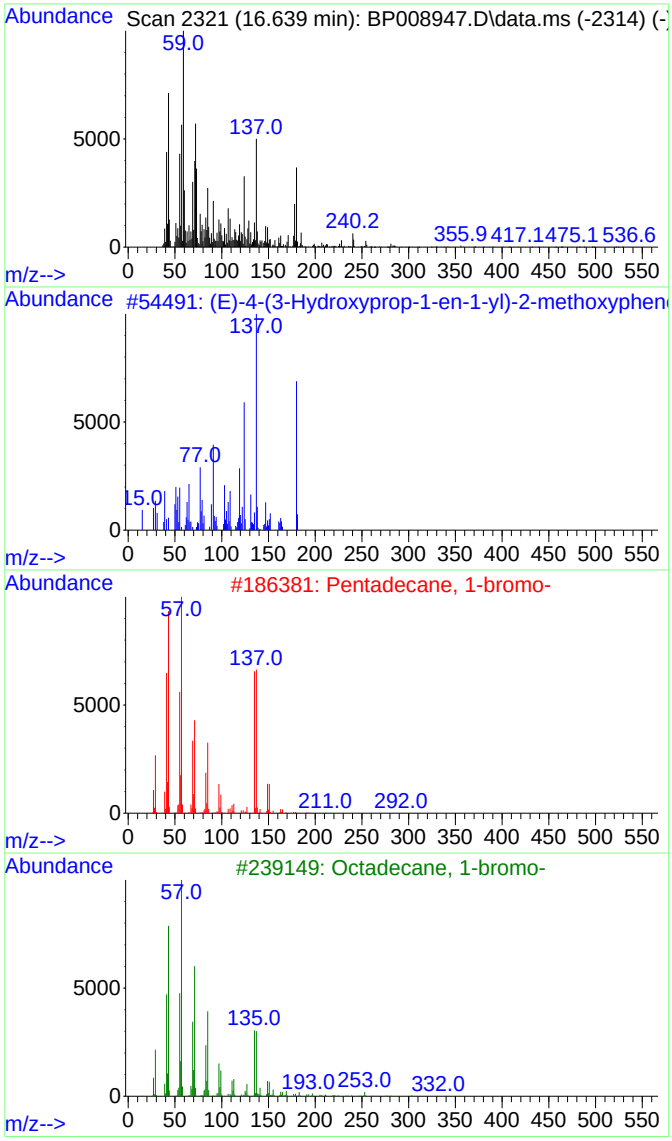
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

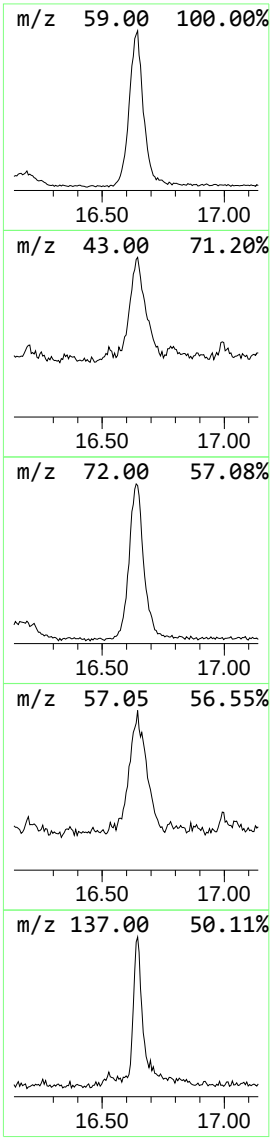
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (E)-4-(3-Hydroxyprop-1-en-1... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	6.51 ng	933616	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	59
2	Pentadecane, 1-bromo-	290	C15H31Br	000629-72-1	53
3	Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	43
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	42
5	Decane, 1-bromo-	220	C10H21Br	000112-29-8	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

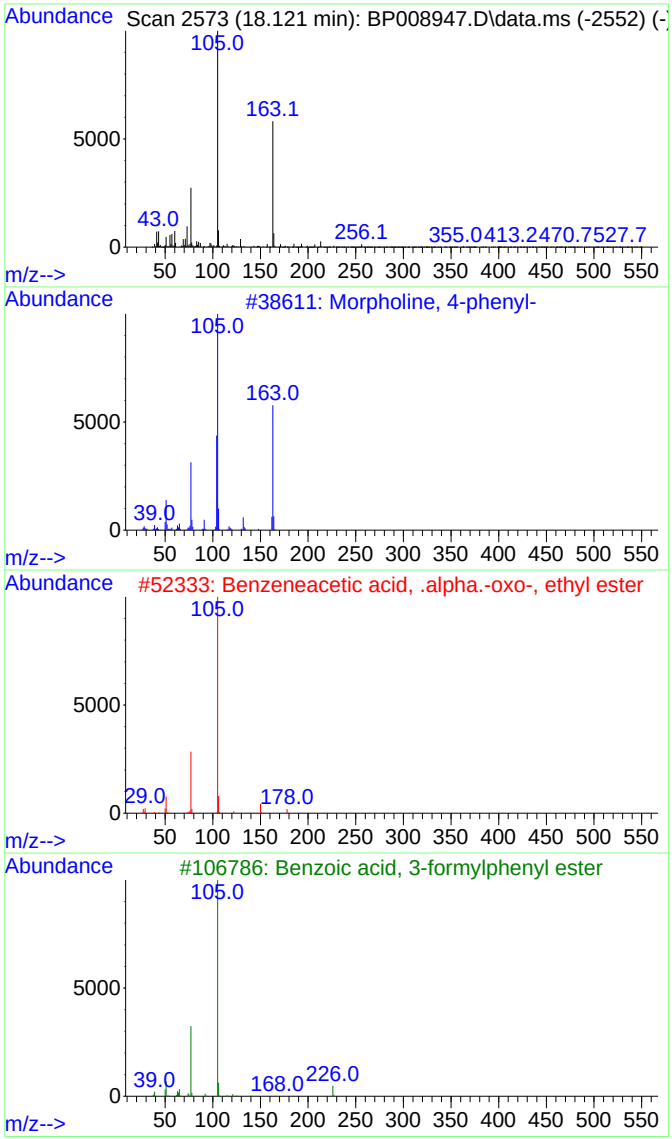
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

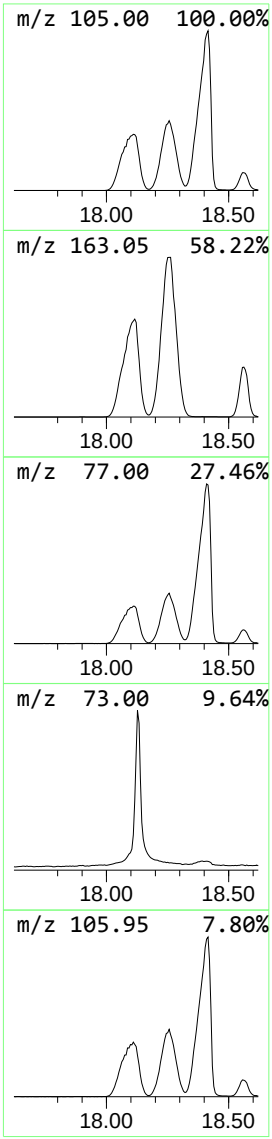
****

Peak Number 3 Morpholine, 4-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.121	228.08 ng	32691600	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	59
2	Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	27
3	Benzoic acid, 3-formylphenyl ester	226	C14H10O3	033696-06-9	27
4	Methanone, (5,6-dimethyl-1H-thie...	256	C14H12N2OS	1000350-07-2	27
5	2,6-Dimethylphenyl benzoate	226	C15H14O2	001871-36-9	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

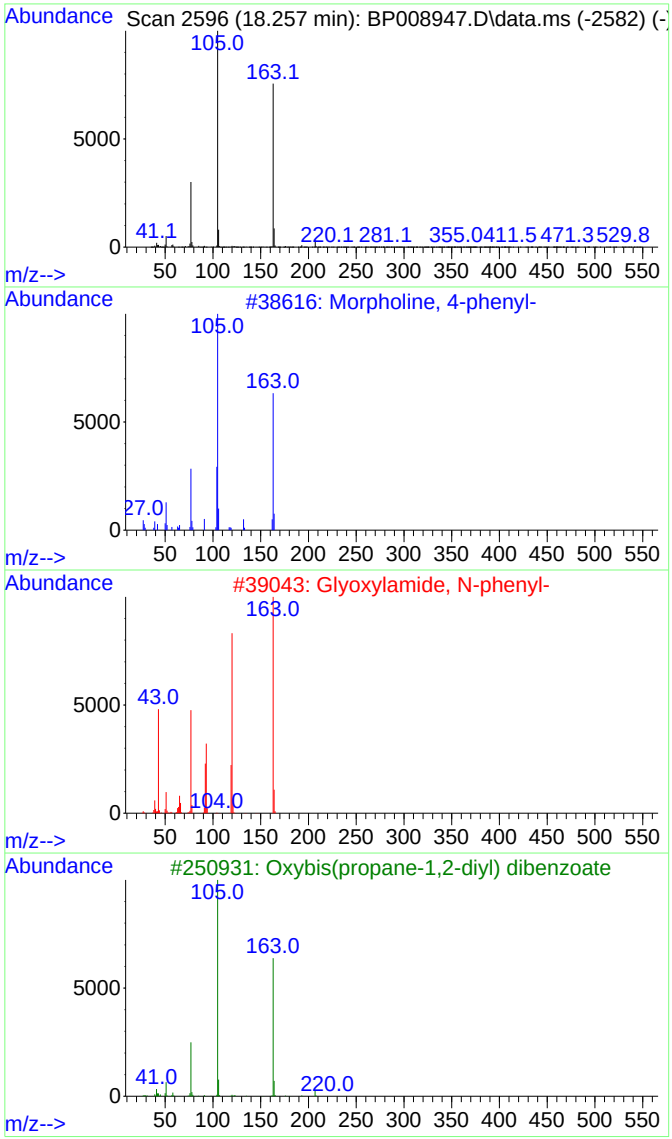
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

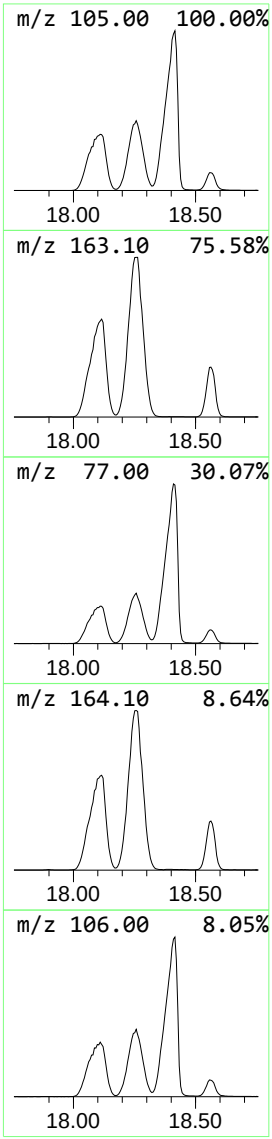
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Glyoxylamide, N-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.257	229.13 ng	32842800	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	78
2	Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	59
3	Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	49
4	4-Nitrophenyl benzoate	243	C13H9NO4	000959-22-8	22
5	Phenylglyoxylic acid, propyl ester	192	C11H12O3	1000453-46-7	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

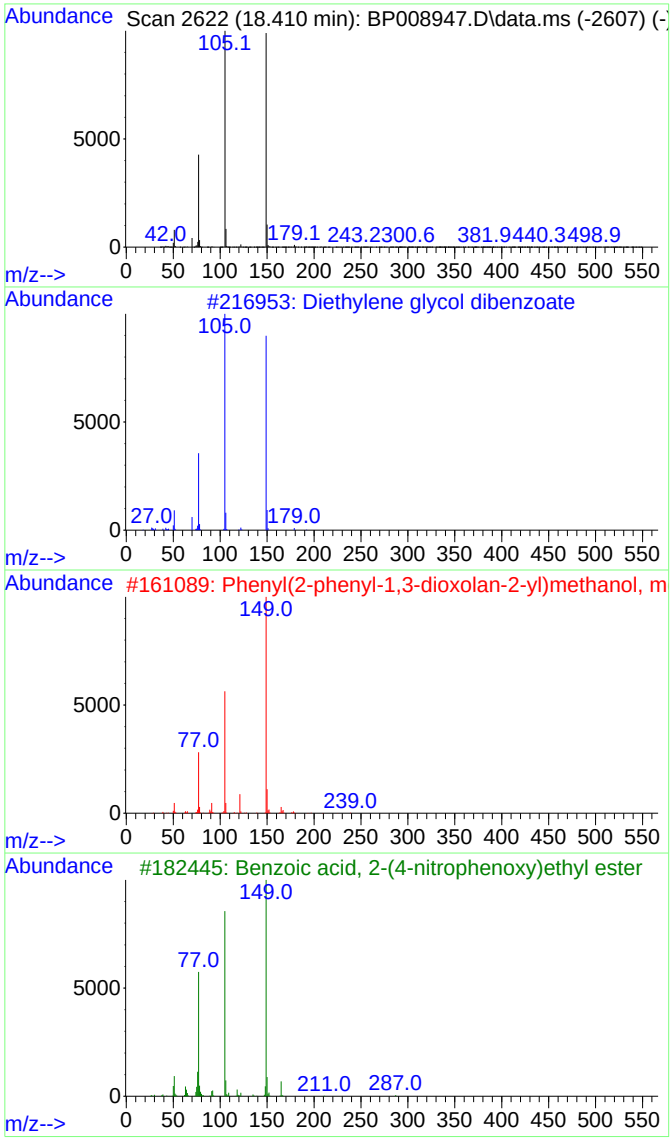
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

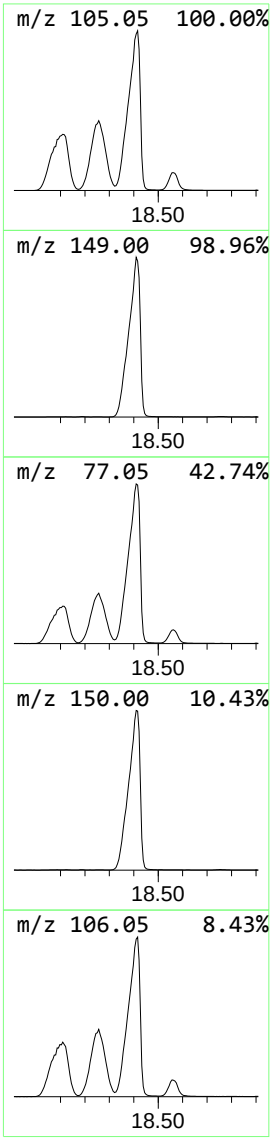
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.410	466.17 ng	66818600	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2	Phenyl(2-phenyl-1,3-dioxolan-2-y...	270	C17H18O3	1000507-07-6	83
3	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	78
4	Phenyl(2-phenyl-1,3-dioxolan-2-y...	298	C18H18O4	006963-16-2	78
5	Phenyl(2-phenyl-1,3-dioxolan-2-y...	256	C16H16O3	005694-69-9	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
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Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

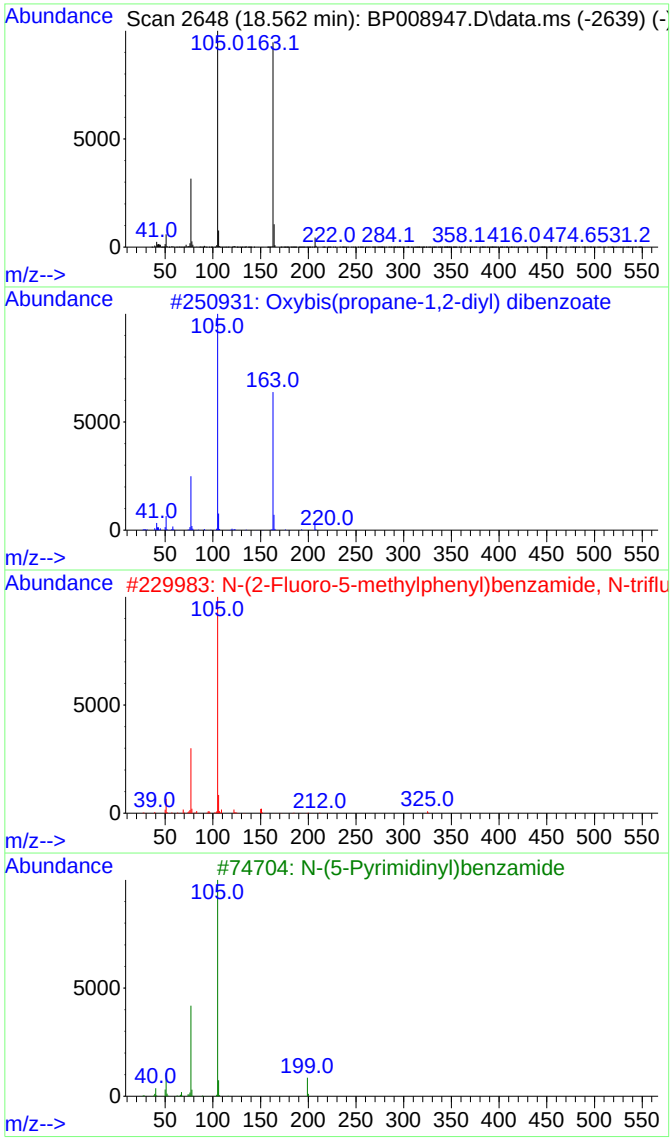
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

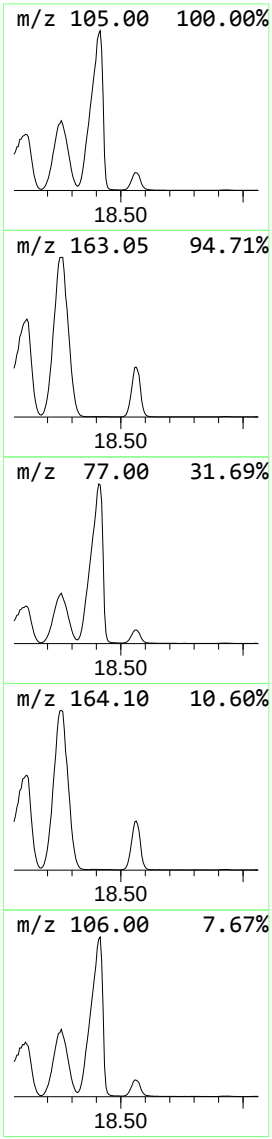
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Oxybis(propane-1,2-diyl) di... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.563	38.31 ng	5491670	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	58
2	N-(2-Fluoro-5-methylphenyl)benza...	325	C16H11F4NO2	1000492-40-0	22
3	N-(5-Pyrimidinyl)benzamide	199	C11H9N3O	103854-65-5	22
4	Phenylglyoxylic acid, isopropyl ...	192	C11H12O3	1000453-46-3	22
5	Benzo[b]thiophene-3-carboxylic a...	343	C18H17NO4S	096334-43-9	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
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Operator : CG/JU
Sample : N1210-02
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
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TIC Library : C:\Database\NIST20.L
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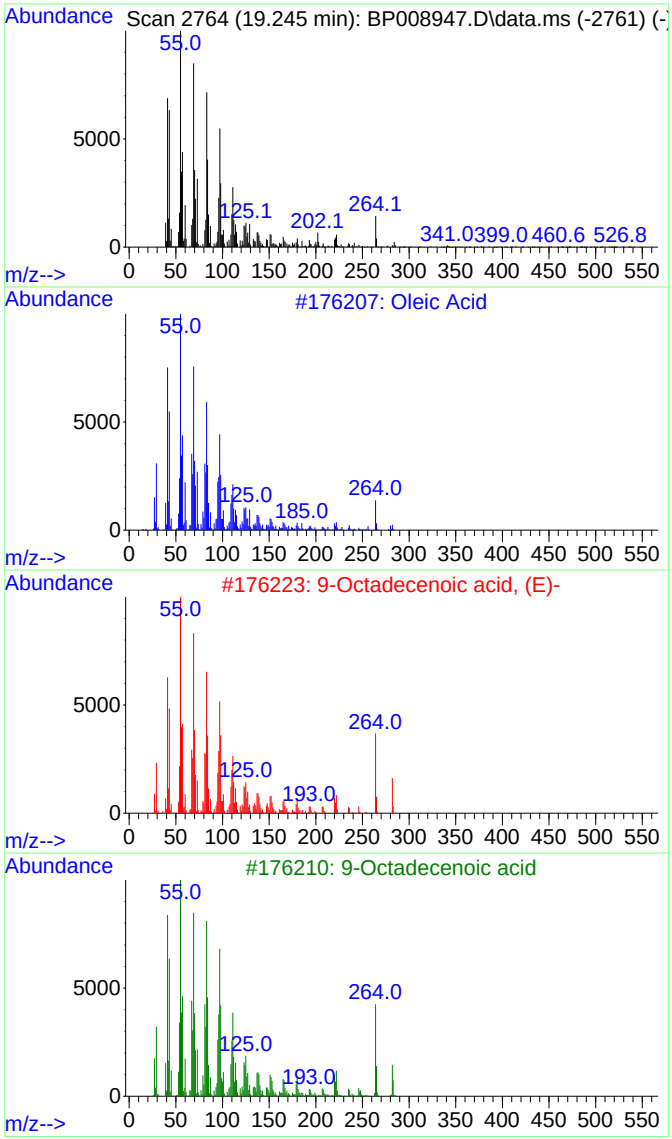
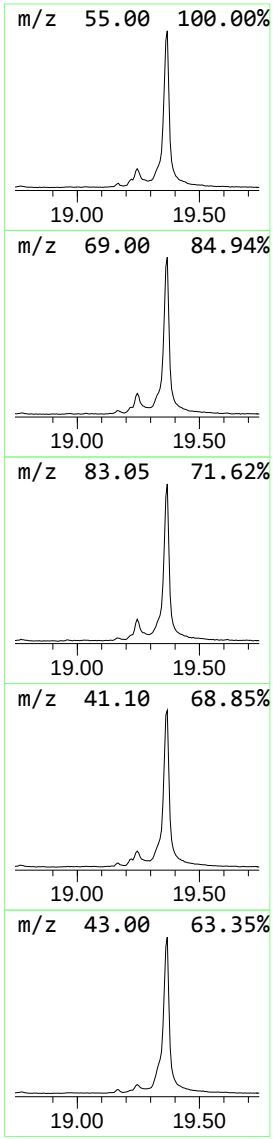
****

Peak Number 7 Oleic Acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	9.00 ng	1289620	Phenanthrene-d10	17.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid	282	C18H34O2	000112-80-1	93
2	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	93
3	9-Octadecenoic acid	282	C18H34O2	002027-47-6	91
4	Oleyl alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	76
5	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	58





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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

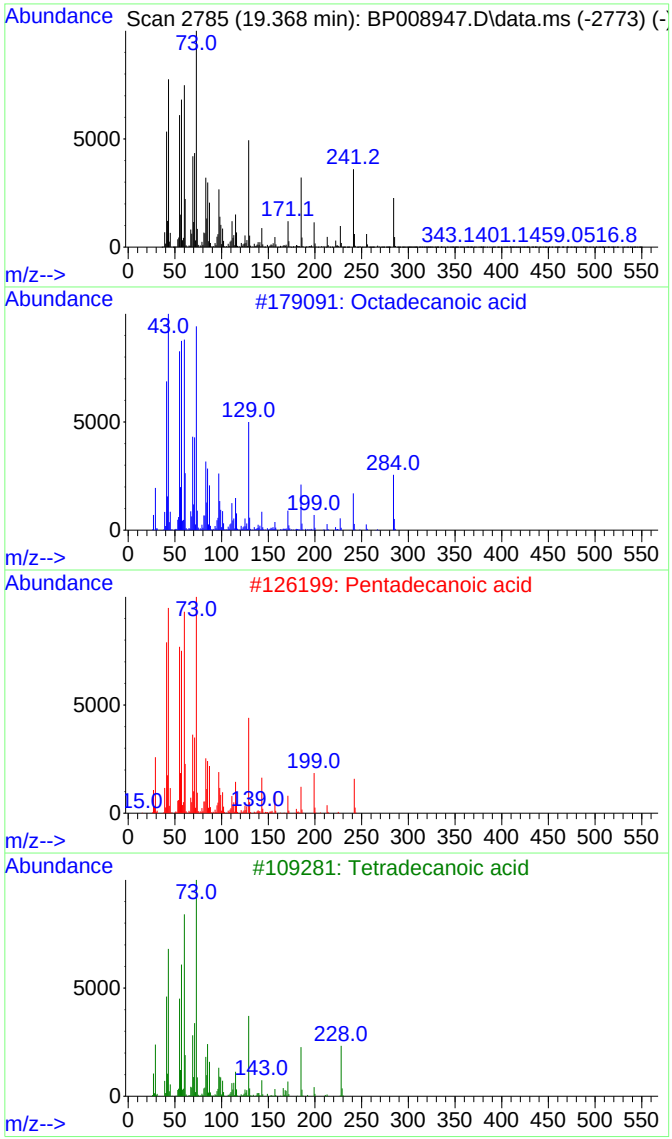
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

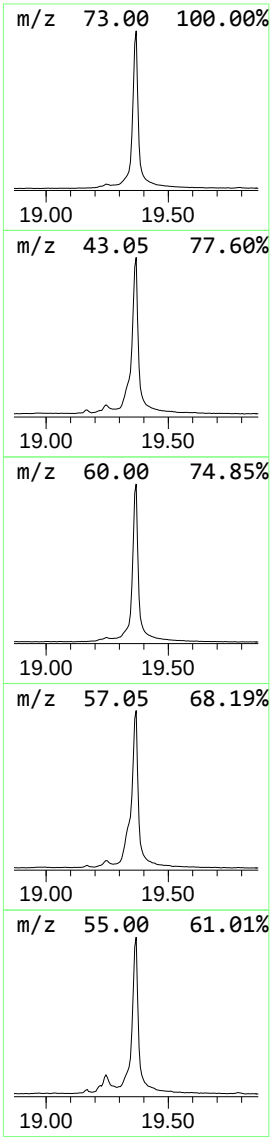
****

Peak Number 8 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.368	85.26 ng	14953400	Chrysene-d12	21.321

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4	Tridecanoic acid	214	C13H26O2	000638-53-9	70
5	n-Decanoic acid	172	C10H20O2	000334-48-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

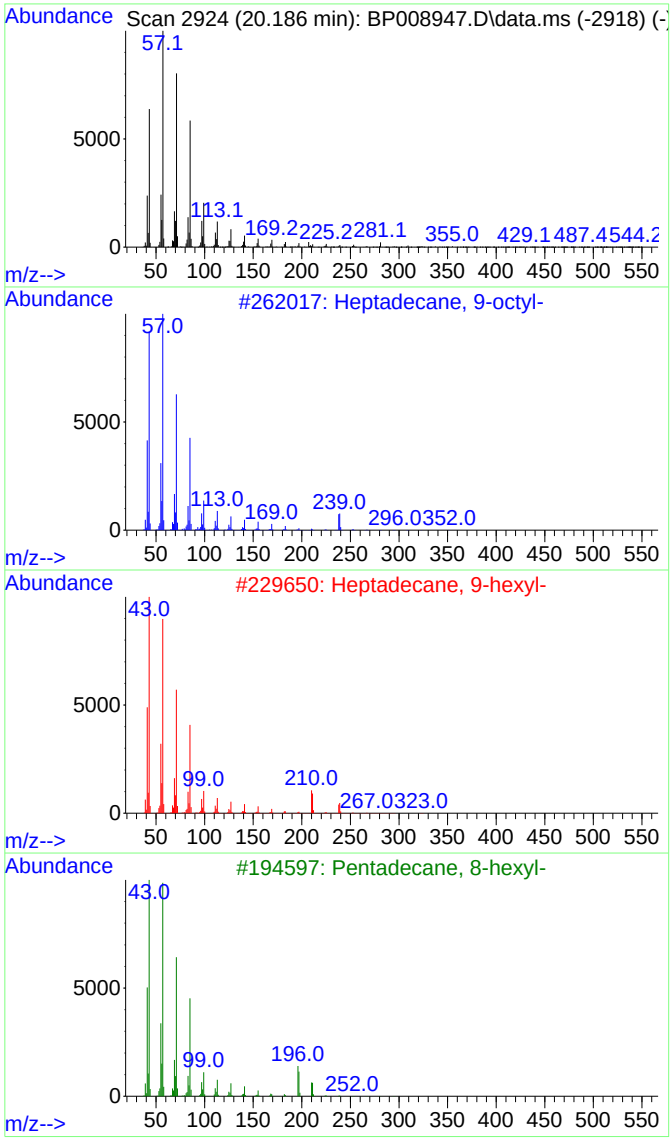
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

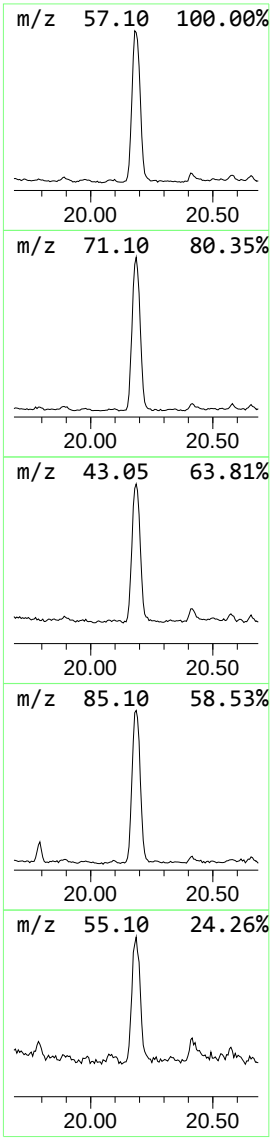
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Heptadecane, 9-octyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.186	7.49 ng	1313510	Chrysene-d12	21.321

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Heneicosane	296	C21H44	000629-94-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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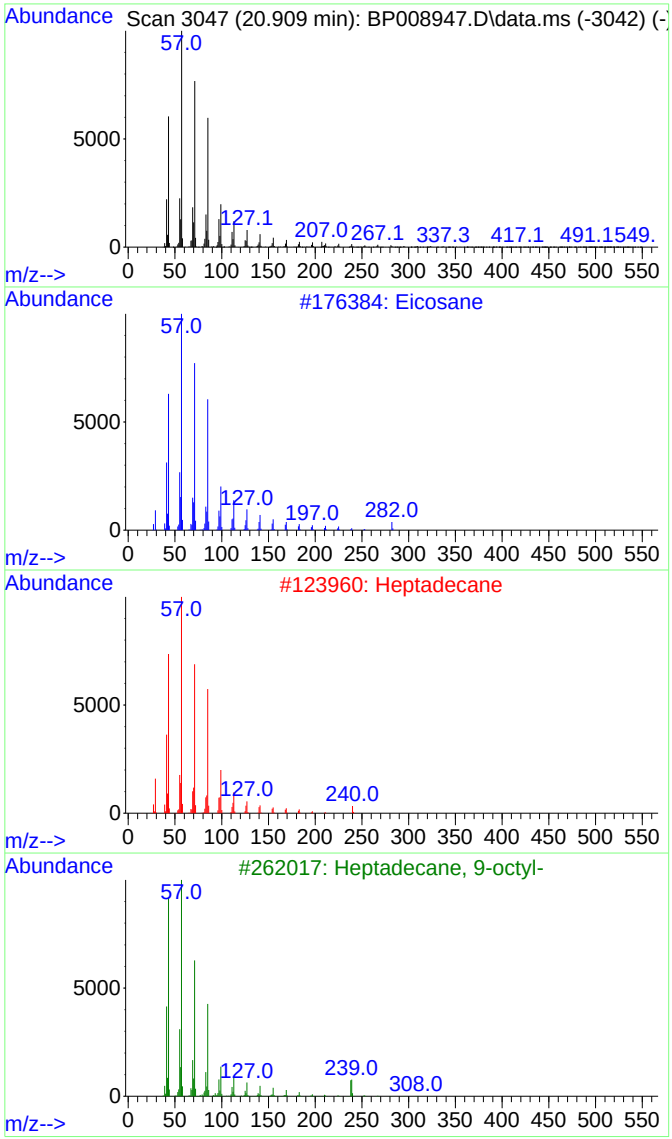
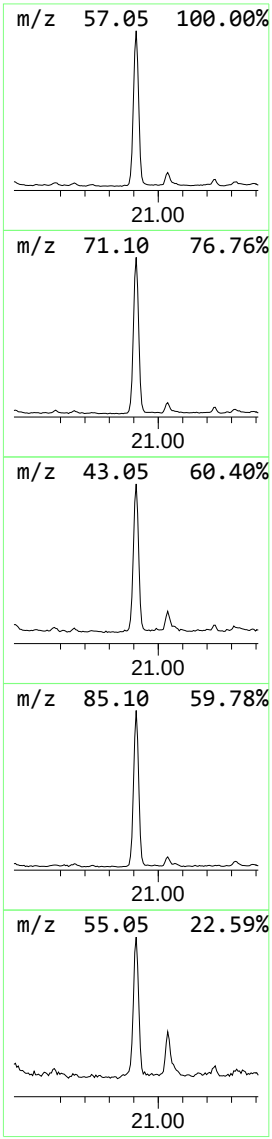
Peak Number 10 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.909	9.87 ng	1730560	Chrysene-d12	21.321

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Heptadecane	240	C17H36	000629-78-7	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
5	Heptacosane	380	C27H56	000593-49-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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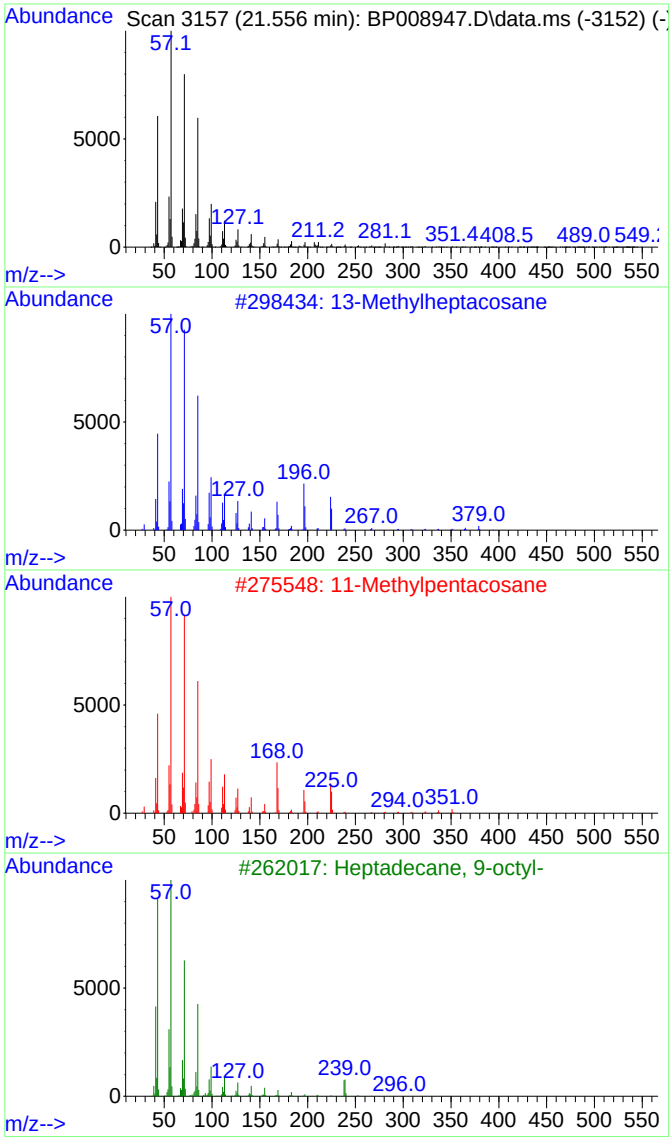
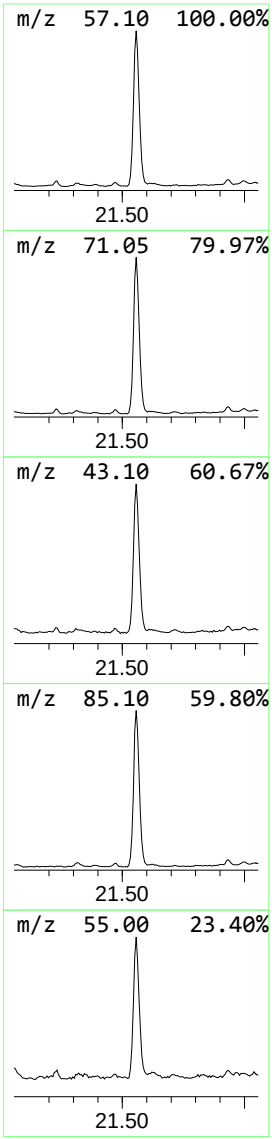
Peak Number 11 13-Methylheptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.556	13.25 ng	2324260	Chrysene-d12	21.321

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Methylheptacosane	394	C28H58	015689-72-2	94
2	11-Methylpentacosane	366	C26H54	015689-71-1	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4	Tetracosane	338	C24H50	000646-31-1	91
5	Heptacosane	380	C27H56	000593-49-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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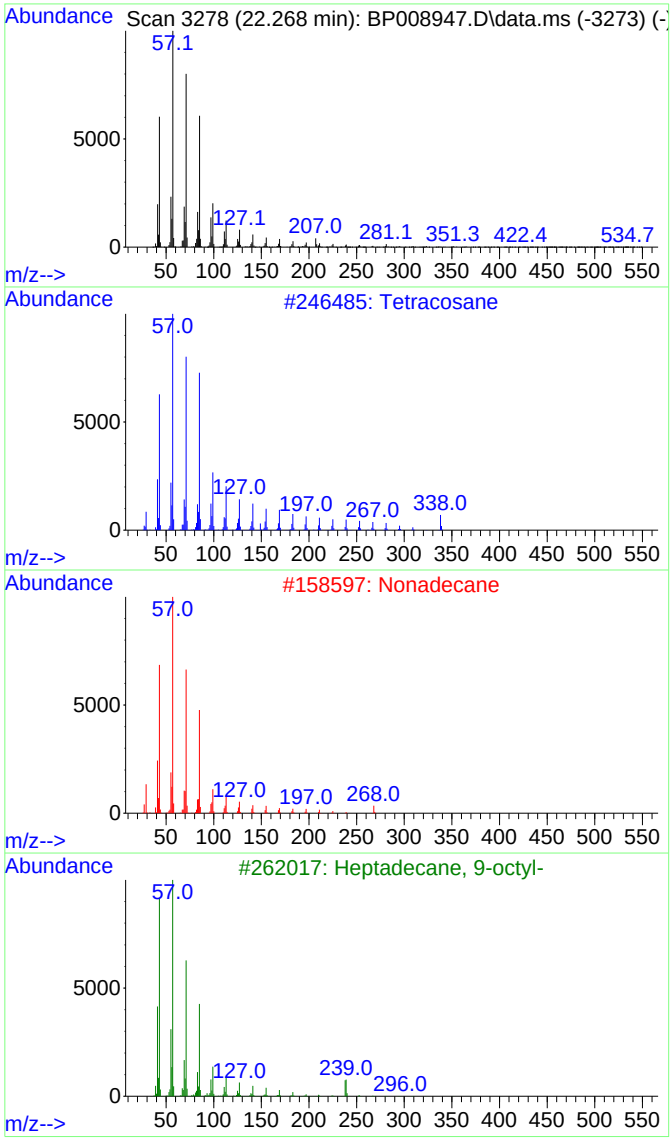
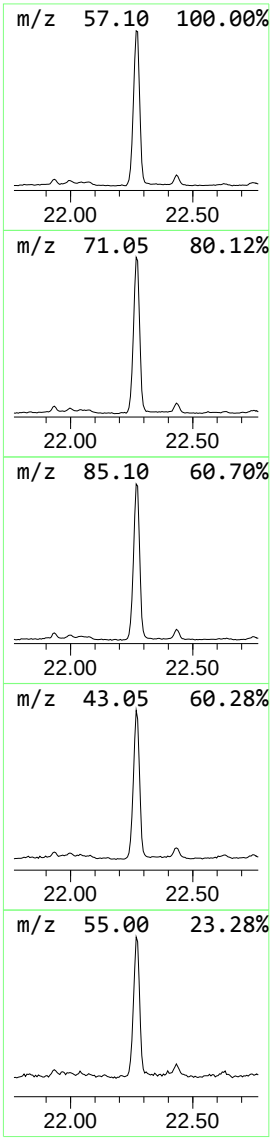
Peak Number 12 Tetracosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.268	13.23 ng	2321200	Chrysene-d12	21.321

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Nonadecane	268	C19H40	000629-92-5	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4	13-Methylheptacosane	394	C28H58	015689-72-2	93
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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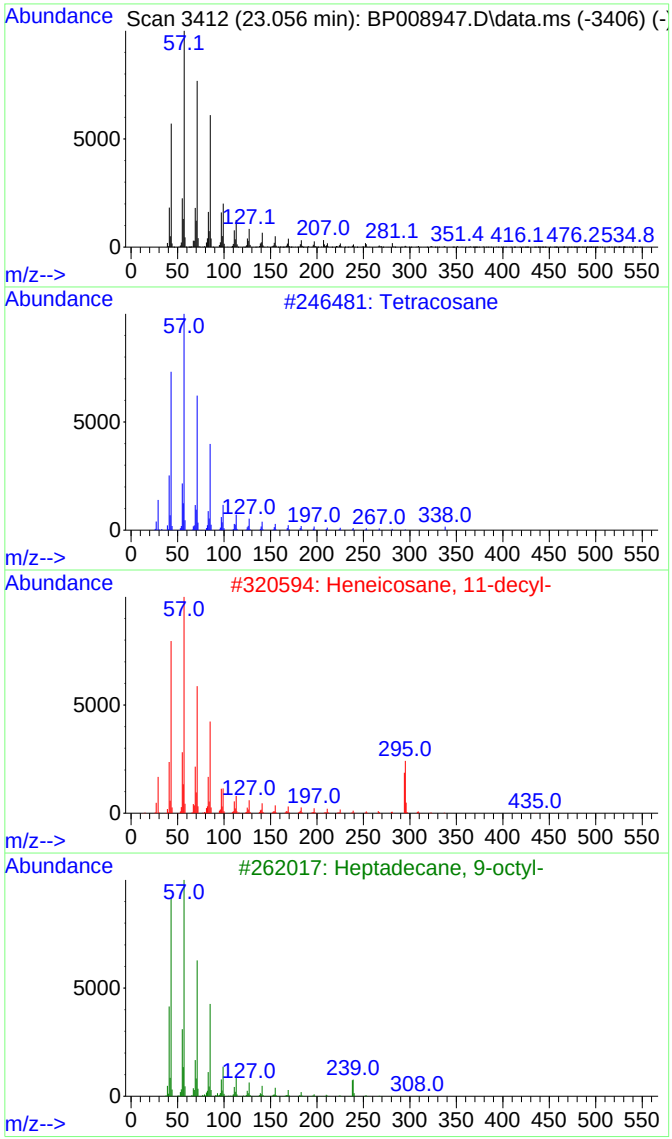
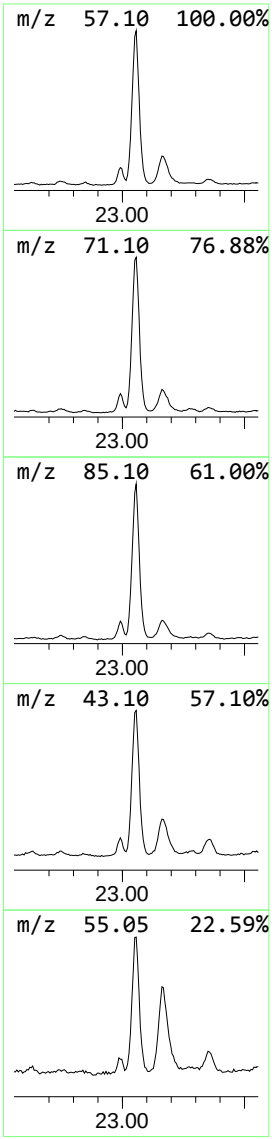
Peak Number 13 Heneicosane, 11-decyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.056	12.85 ng	2315960	Perylene-d12	23.727

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4	11-Methylpentacosane	366	C26H54	015689-71-1	91
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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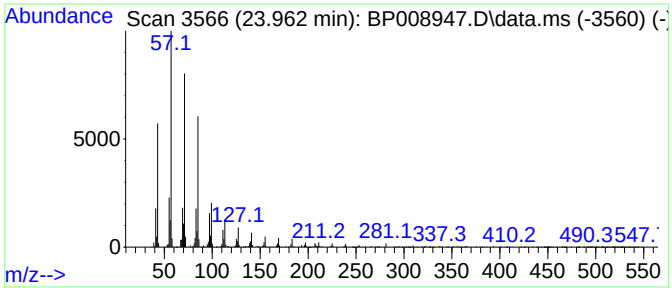
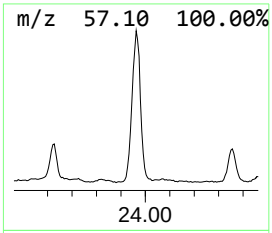
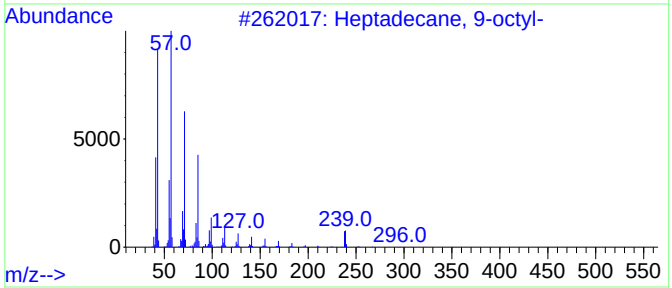
Peak Number 14 Octacosane, 2-methyl- Concentration Rank 11

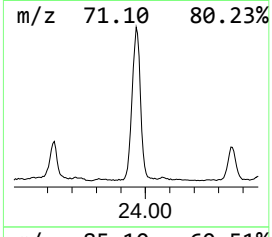
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.962	9.26 ng	1668970	Perylene-d12	23.727

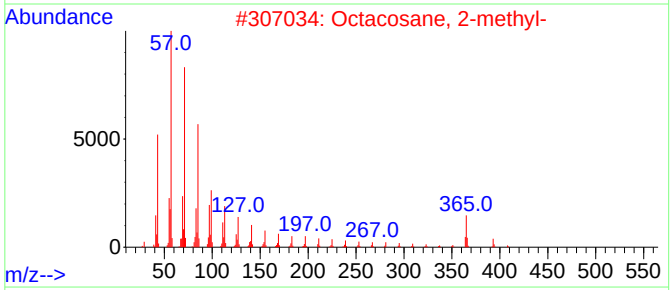
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	94
3	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
4	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
5	2-Methylhentriacontane	451	C32H66	001720-12-3	91

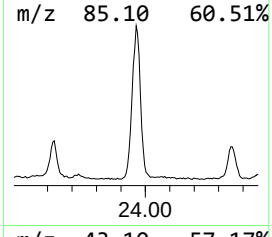


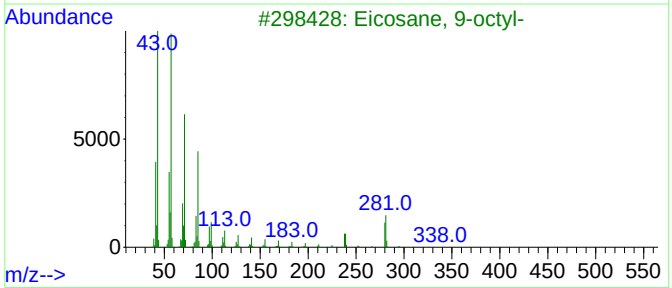


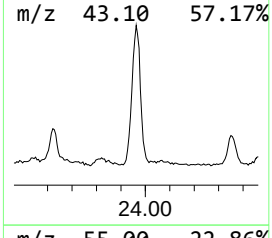


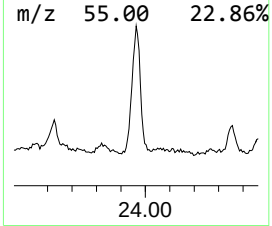












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
Data File : BP008947.D
Acq On : 27 Jan 2022 15:02
Operator : CG/JU
Sample : N1210-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	3.040	32.1	ng	2099820	1	7.828	1307710	20.0
(E)-4-(3-Hydrox...	16.639	6.5	ng	933616	4	17.227	2866730	20.0
Morpholine, 4-p...	18.121	228.1	ng	32691600	4	17.227	2866730	20.0
Glyoxylamide, N...	18.257	229.1	ng	32842800	4	17.227	2866730	20.0
Diethylene glyc...	18.410	466.2	ng	66818600	4	17.227	2866730	20.0
Oxybis(propane-...	18.563	38.3	ng	5491670	4	17.227	2866730	20.0
Oleic Acid	19.245	9.0	ng	1289620	4	17.227	2866730	20.0
Octadecanoic acid	19.368	85.3	ng	14953400	5	21.321	3507780	20.0
Heptadecane, 9-...	20.186	7.5	ng	1313510	5	21.321	3507780	20.0
Eicosane	20.909	9.9	ng	1730560	5	21.321	3507780	20.0
13-Methylheptac...	21.556	13.3	ng	2324260	5	21.321	3507780	20.0
Tetracosane	22.268	13.2	ng	2321200	5	21.321	3507780	20.0
Heneicosane, 11...	23.056	12.9	ng	2315960	6	23.727	3604270	20.0
Octacosane, 2-m...	23.962	9.3	ng	1668970	6	23.727	3604270	20.0