LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051021\
Data File : BP005464.D
Acq On : 10 May 2021 12:39
Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

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-BP050521.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005464.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.828	288	296	303	rVB2	8136	13303	1.13%	0.248%
2	5.293	368	375	387	rBV	253980	347310	29.62%	6.483%
3	6.899	641	648	663	rBV	197889	320535	27.34%	5.983%
4	7.699	778	784	791	rBV	52468	82190	7.01%	1.534%
5	8.869	975	983	997	rBV	135385	218612	18.64%	4.081%
6	10.493	1252	1259	1267	rBV	51277	86514	7.38%	1.615%
7	12.975	1674	1681	1690	rBV	355120	514881	43.91%	9.611%
8	13.822	1820	1825	1832	rBV	9391	15337	1.31%	0.286%
9	14.351	1909	1915	1924	rVB	81069	117225	10.00%	2.188%
10	15.845	2161	2169	2181	rBV	404660	567998	48.44%	10.602%
11	16.516	2278	2283	2295	rVB6	25010	47054	4.01%	0.878%
12	17.104	2377	2383	2390	rBV	118298	161869	13.81%	3.021%
13	17.775	2485	2497	2507	rBV9	20224	64850	5.53%	1.211%
14	17.928	2512	2523	2530	rBV2	25453	79034	6.74%	1.475%
15	18.004	2531	2536	2544	rBV	96318	136181	11.61%	2.542%
16	18.280	2577	2583	2587	rBV2	50533	78871	6.73%	1.472%
17	18.327	2587	2591	2613	rVB5	77492	207637	17.71%	3.876%
18	19.027	2706	2710	2719	rBV8	10989	24627	2.10%	0.460%
19	19.127	2724	2727	2734	rVB8	10427	16770	1.43%	0.313%
20	19.239	2742	2746	2760	rVB2	62395	85734	7.31%	1.600%
21	19.674	2814	2820	2827	rBV	1061002	1172518	100.00%	21.886%
22	19.898	2853	2858	2861	rBV7	7689	12826	1.09%	0.239%
23	19.963	2864	2869	2875	rVB7	16481	37294	3.18%	0.696%
24	20.374	2933	2939	2945	rBV2	38932	69699	5.94%	1.301%
25	20.721	2994	2998	3005	rVB5	20988	38366	3.27%	0.716%
26	20.880	3019	3025	3027	rBV3	26869	48082	4.10%	0.898%
27	20.910	3027	3030	3037	rVV4	54037	104419	8.91%	1.949%
28	20.980	3039	3042	3048	rVB	49580	61121	5.21%	1.141%
29	21.198	3074	3079	3089	rBV	189667	251568	21.46%	4.696%
30	21.386	3107	3111	3116	rVB2	34650	50871	4.34%	0.950%
31	23.515	3466	3473	3484	rVB2	154257	323983	27.63%	6.048%

Sum of corrected areas: 5357279

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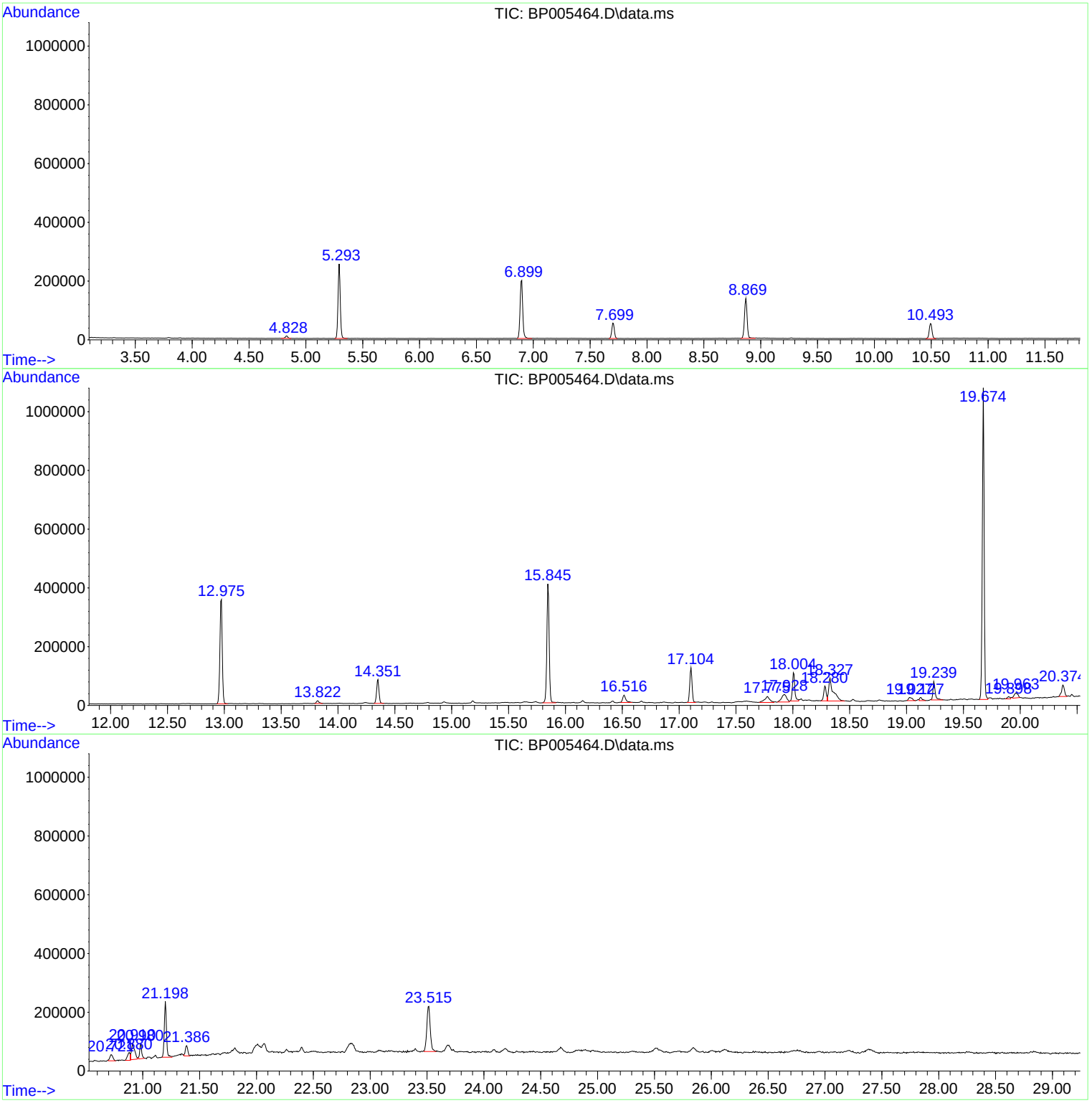
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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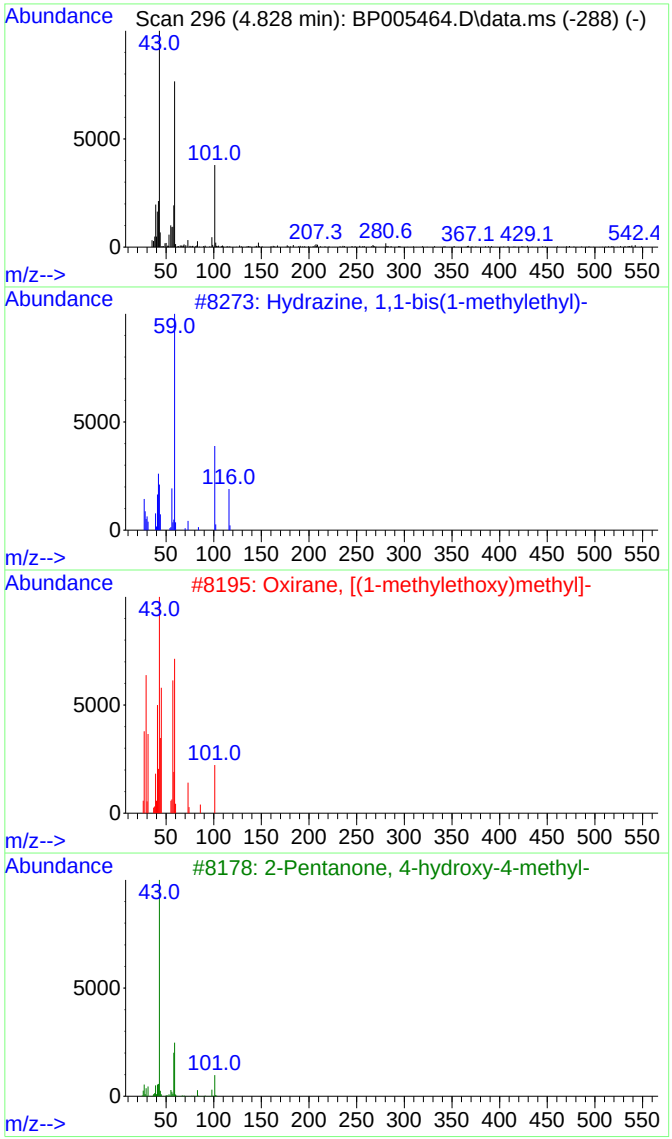
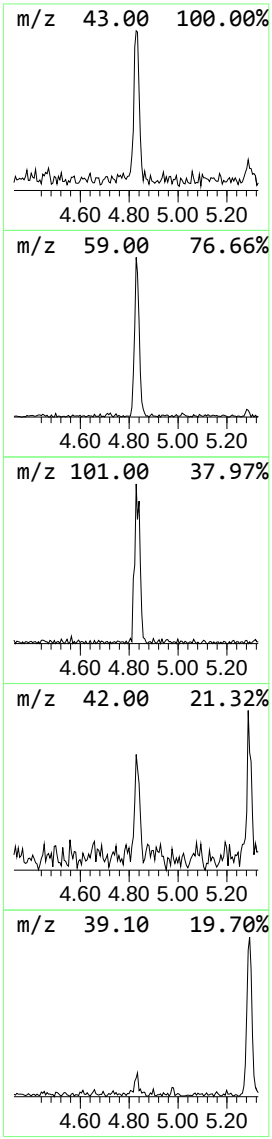
****
Peak Number 1 unknown4.82 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.828	3.24 ng	13303	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	39
2	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	39
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39
4	Hexanamide	115	C6H13NO	000628-02-4	33
5	2-Pentanol, 5-methoxy-2-methyl-	132	C7H16O2	055724-04-4	33





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Data File : BP005464.D
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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

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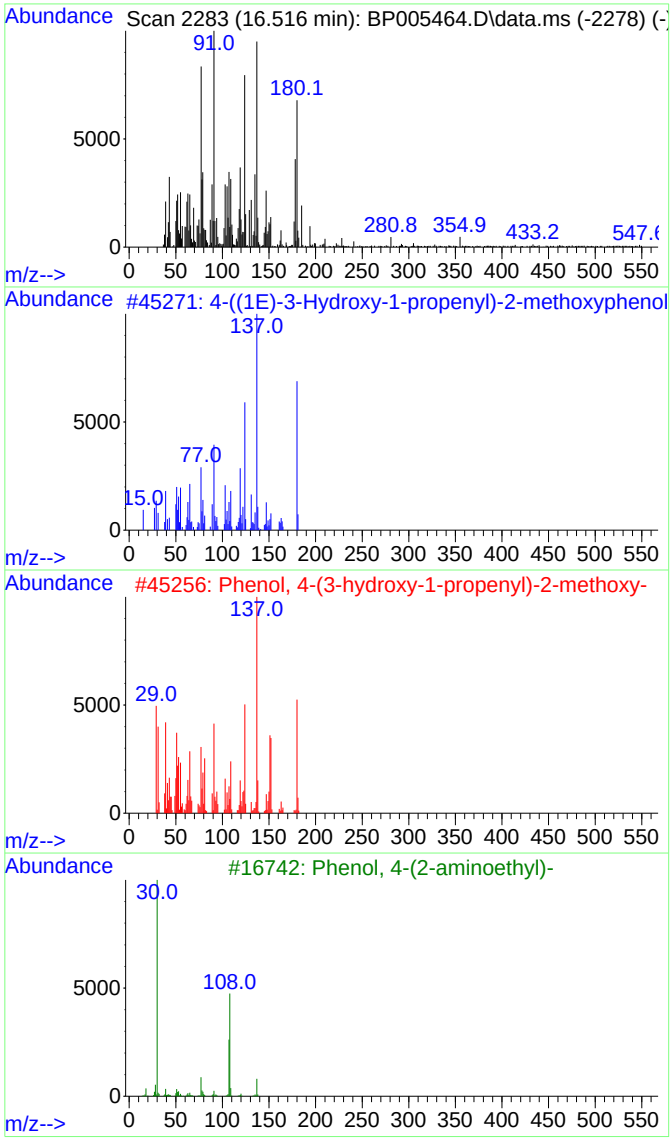
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

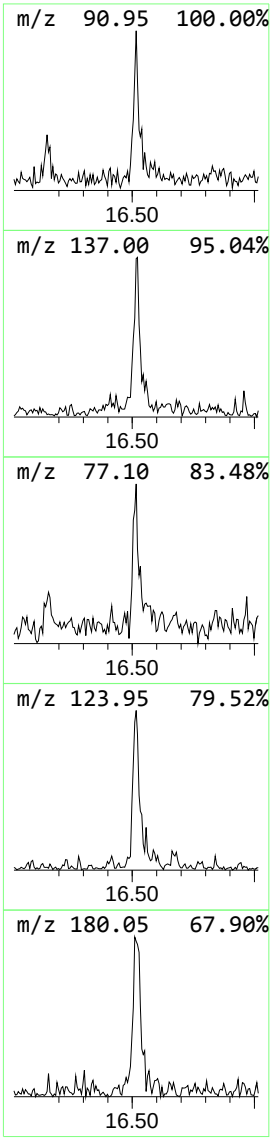
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 4-((1E)-3-Hydroxy-1-propeny... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.516	5.81 ng	47054	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-((1E)-3-Hydroxy-1-propenyl)-2-...	180	C10H12O3	1000297-95-5	56
2	Phenol, 4-(3-hydroxy-1-propenyl)...	180	C10H12O3	000458-35-5	41
3	Phenol, 4-(2-aminoethyl)-	137	C8H11NO	000051-67-2	35
4	9-Ethoxy-10-oxatricyclo[7.2.1.0(...	224	C13H20O3	1000187-02-8	25
5	2-Methylthianaphthene-1,1 dioxide	180	C9H8O2S	006224-55-1	20





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ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

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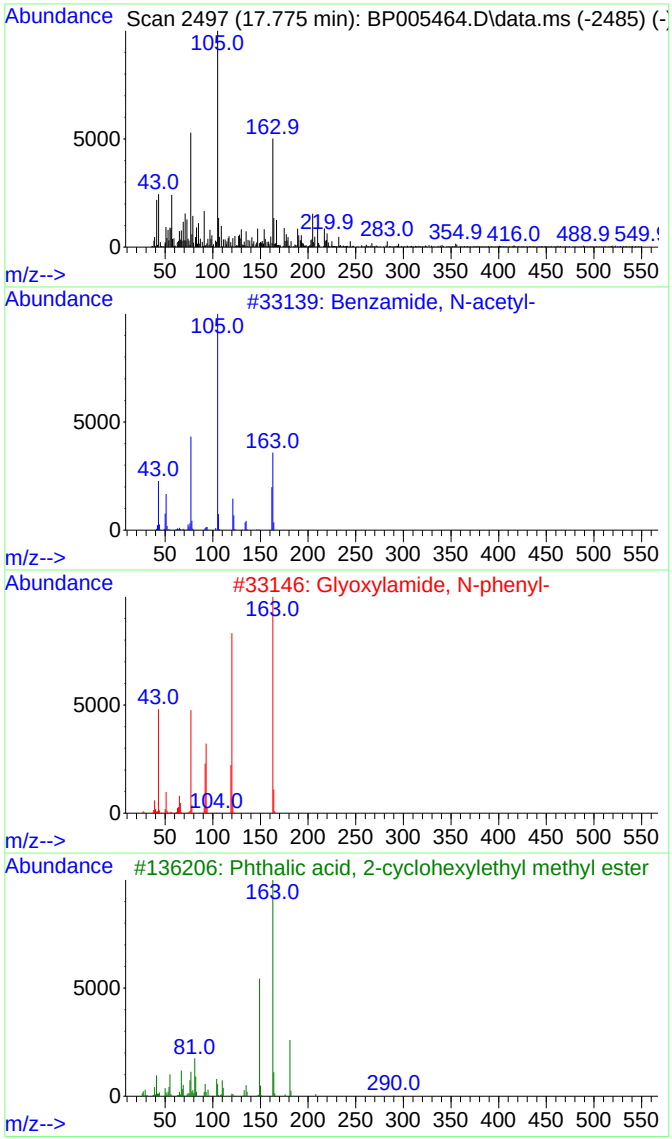
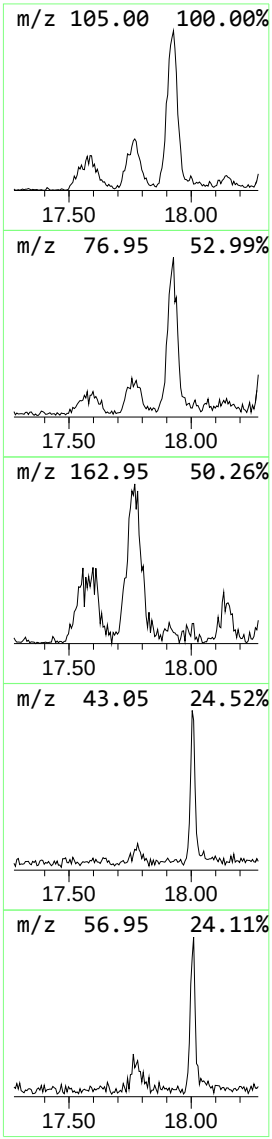
****

Peak Number 3 Benzamide, N-acetyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.775	8.01 ng	64850	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	64
2	Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	47
3	Phthalic acid, 2-cyclohexylethyl...	290	C17H22O4	1000309-03-4	38
4	N,N-Diisopropylbenzamide	205	C13H19NO	020383-28-2	38
5	4H-1,3,2-Dioxaborin, 4,6-diethen...	178	C10H15BO2	074630-03-8	35





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

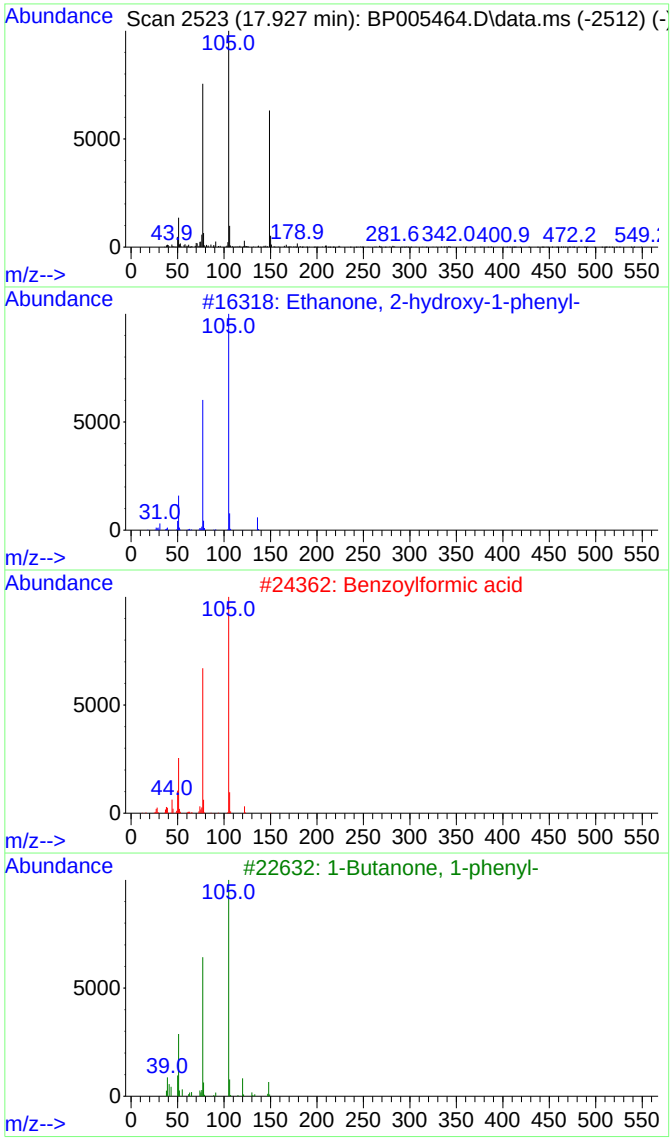
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

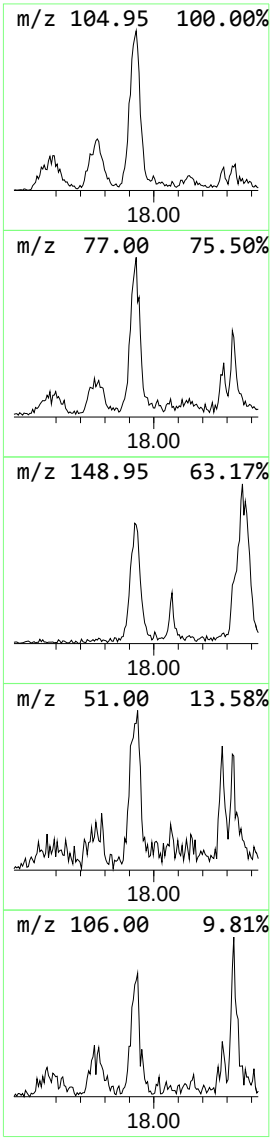
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanone, 2-hydroxy-1-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.927	9.77 ng	79034	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 2-hydroxy-1-phenyl-	136	C8H8O2	000582-24-1	53
2	Benzoylformic acid	150	C8H6O3	000611-73-4	53
3	1-Butanone, 1-phenyl-	148	C10H12O	000495-40-9	53
4	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	53
5	Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	53





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

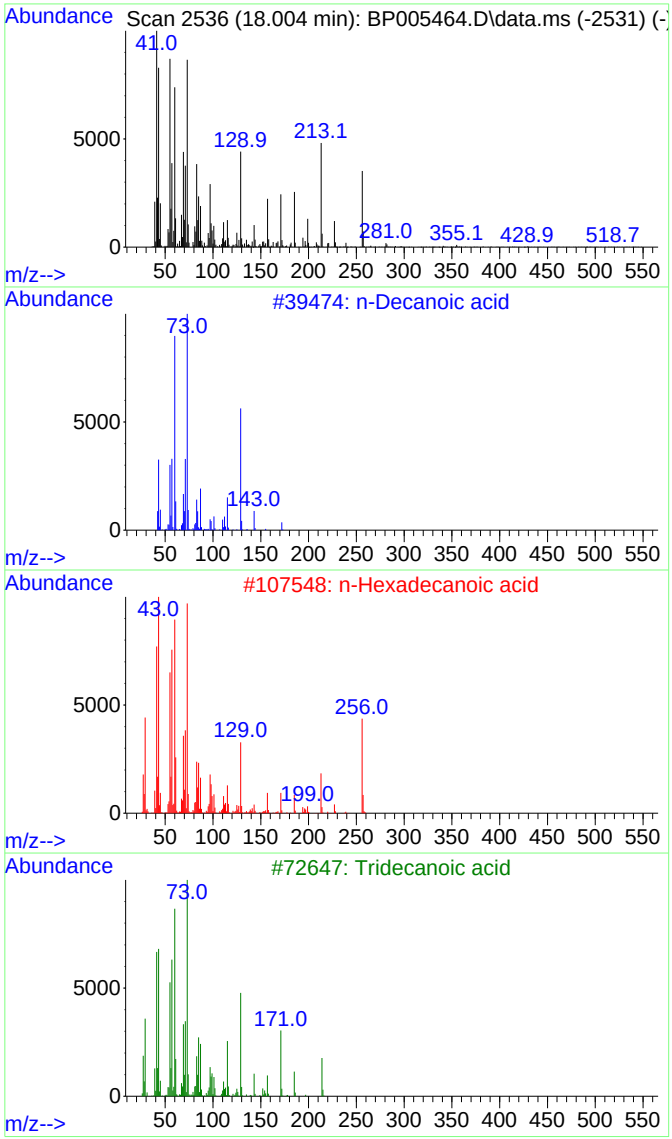
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

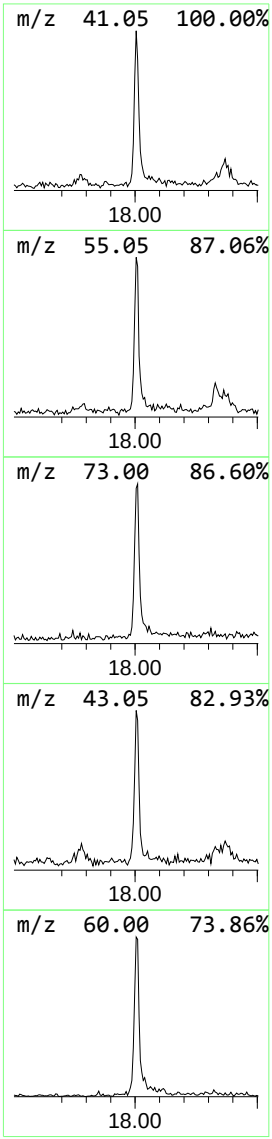
****

Peak Number 5 n-Decanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.004	16.83 ng	136181	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	80
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	78
3	Tridecanoic acid	214	C13H26O2	000638-53-9	46
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	43
5	Octadecanoic acid	284	C18H36O2	000057-11-4	43





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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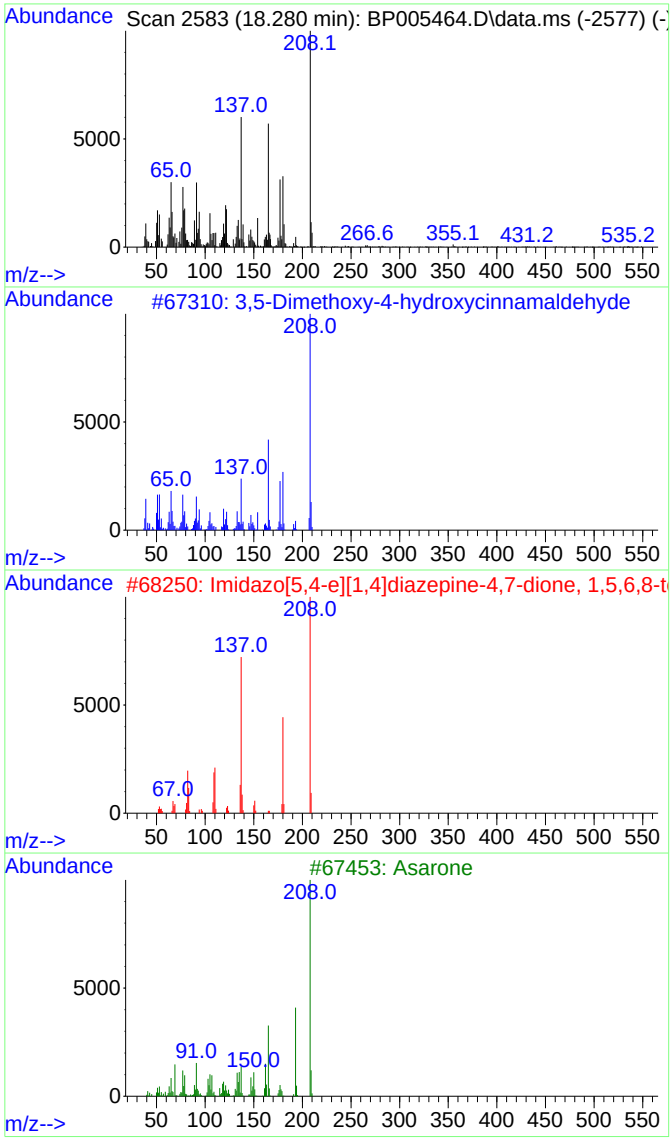
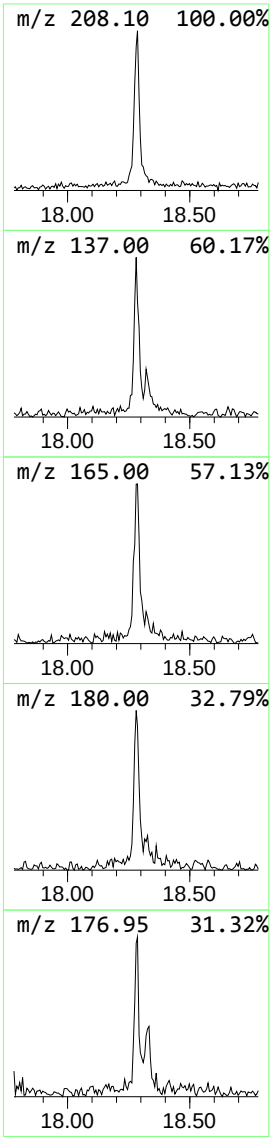
****
Peak Number 6 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.280	9.75 ng	78871	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethoxy-4-hydroxycinnamald...	208	C11H12O4	087345-53-7	90
2	Imidazo[5,4-e][1,4]diazepine-4,7...	208	C9H12N4O2	144105-22-6	50
3	Asarone	208	C12H16O3	002883-98-9	46
4	2-Propenoic acid, 3-(2,3-dimetho...	208	C11H12O4	007345-82-6	41
5	Acetic acid, (4-formylphenoxy)-,...	208	C11H12O4	051264-69-8	38





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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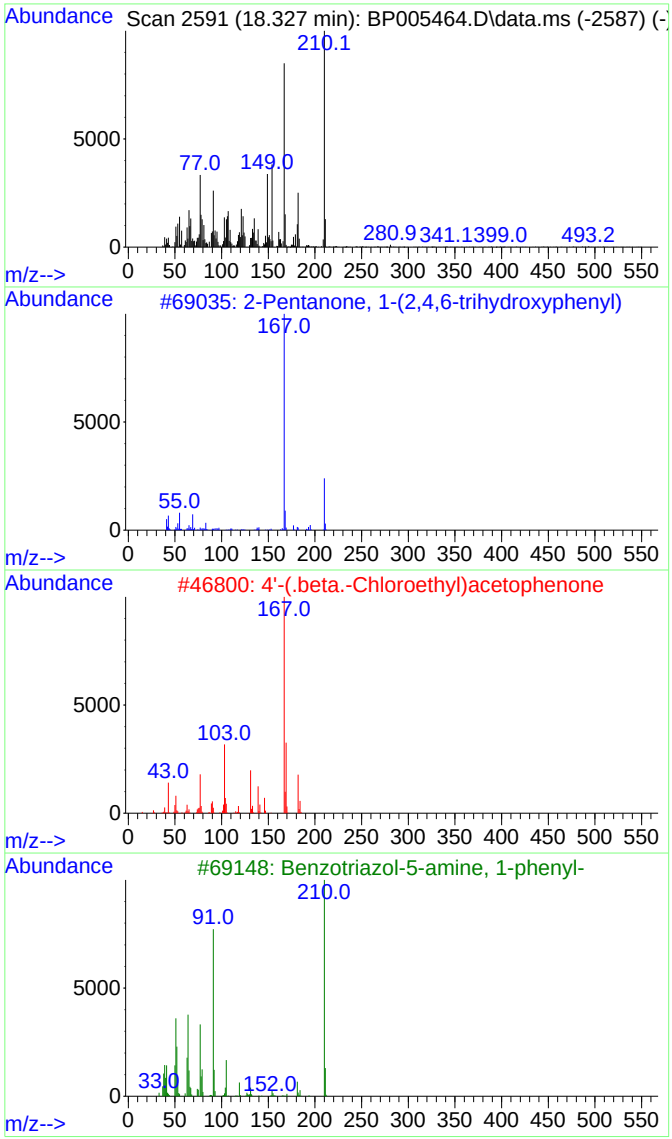
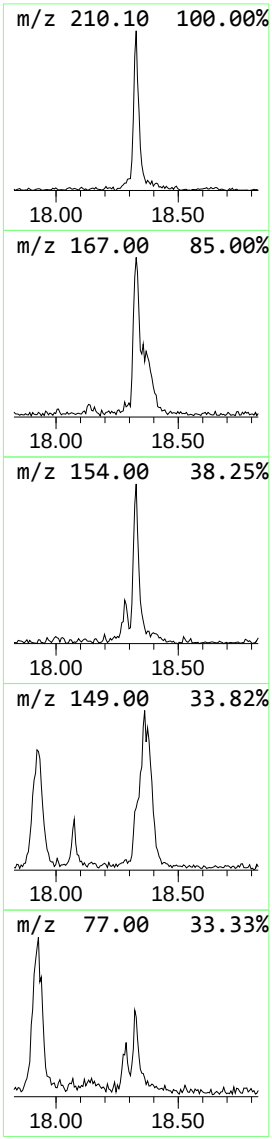
****
Peak Number 7 2-Pentanone, 1-(2,4,6-trihy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.328	25.65 ng	207637	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	50
2	4'-(.beta.-Chloroethyl)acetophenone	182	C10H11ClO	069614-95-5	30
3	Benzotriazol-5-amine, 1-phenyl-	210	C12H10N4	021819-66-9	30
4	Phenol, 2-(2-imidazo[1,2-a]pyrid...	210	C13H10N2O	057636-32-5	27
5	Benzene, 1,1'-ethylidenebis-	182	C14H14	000612-00-0	27





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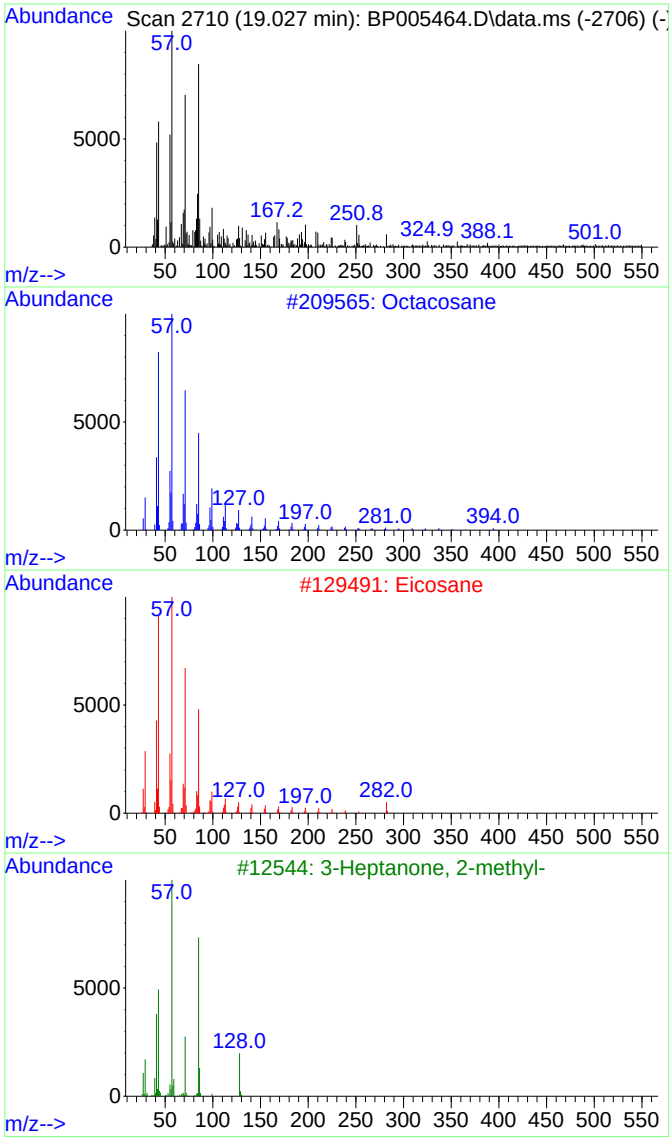
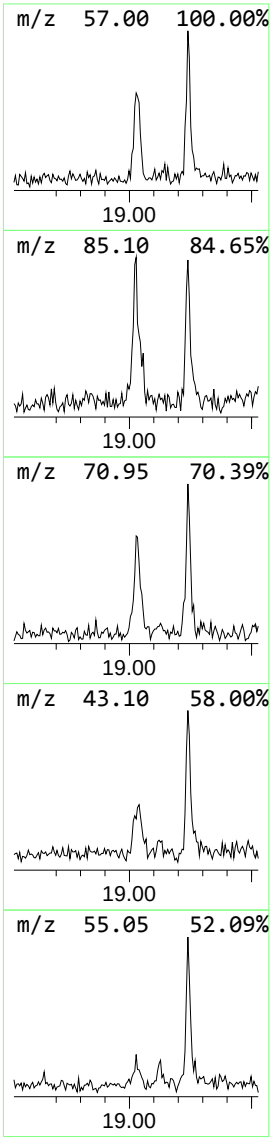
****

Peak Number 8 Octacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.027	3.04 ng	24627	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	64
2	Eicosane	282	C20H42	000112-95-8	64
3	3-Heptanone, 2-methyl-	128	C8H16O	013019-20-0	49
4	Pentacosane	352	C25H52	000629-99-2	43
5	2-Bromotetradecane	276	C14H29Br	074036-95-6	43





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Sample : M2306-01
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ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
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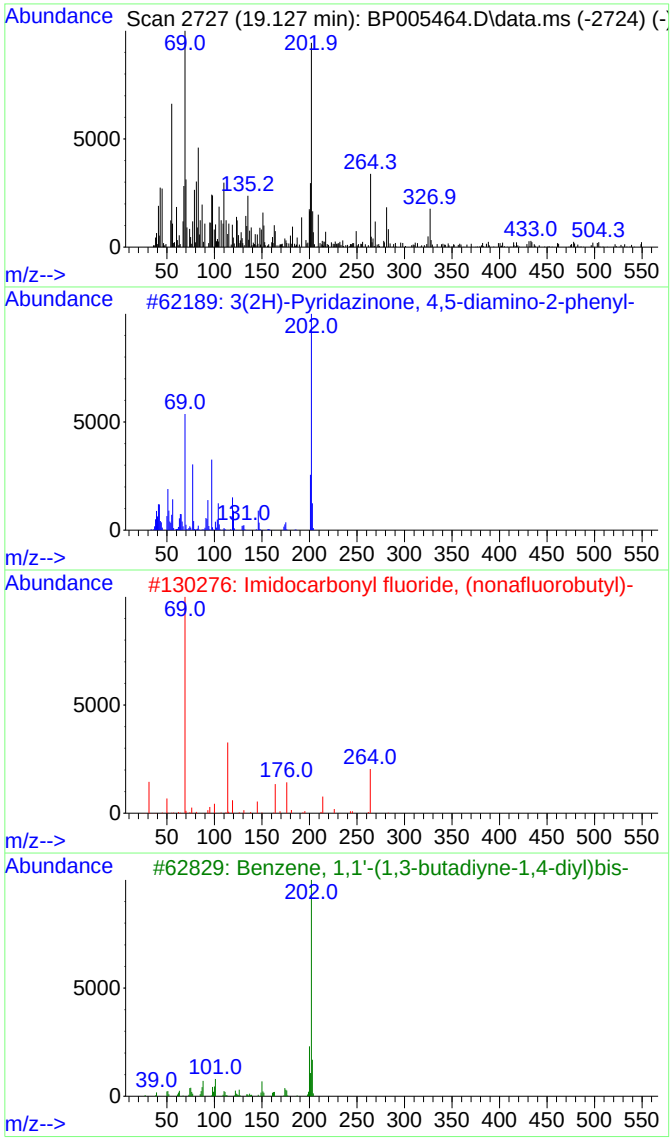
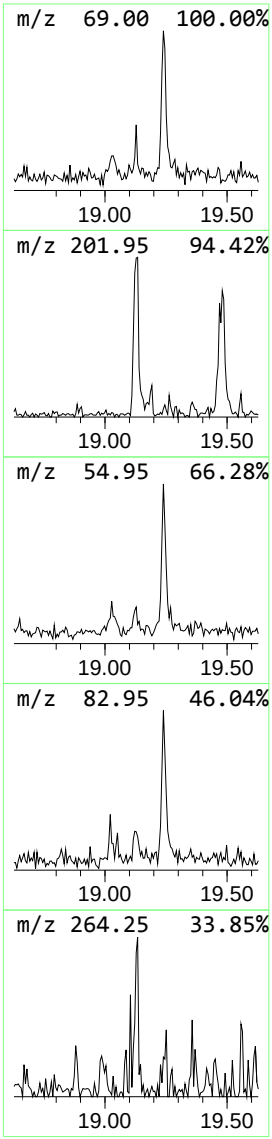
****
Peak Number 9 3(2H)-Pyridazinone, 4,5-dia... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.127	2.07 ng	16770	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3(2H)-Pyridazinone, 4,5-diamino-...	202	C10H10N4O	1000351-65-9	64
2	Imidocarbonyl fluoride, (nonaflu...	283	C5F11N	000424-32-8	38
3	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	30
4	Cyclopropane-1,1-dicarbonitrile,...	202	C11H7ClN2	098235-21-3	30
5	Phenol, 2-(5-bromo-2-thienyliden...	281	C11H8BrNOS	099303-13-6	30





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Acq On : 10 May 2021 12:39
Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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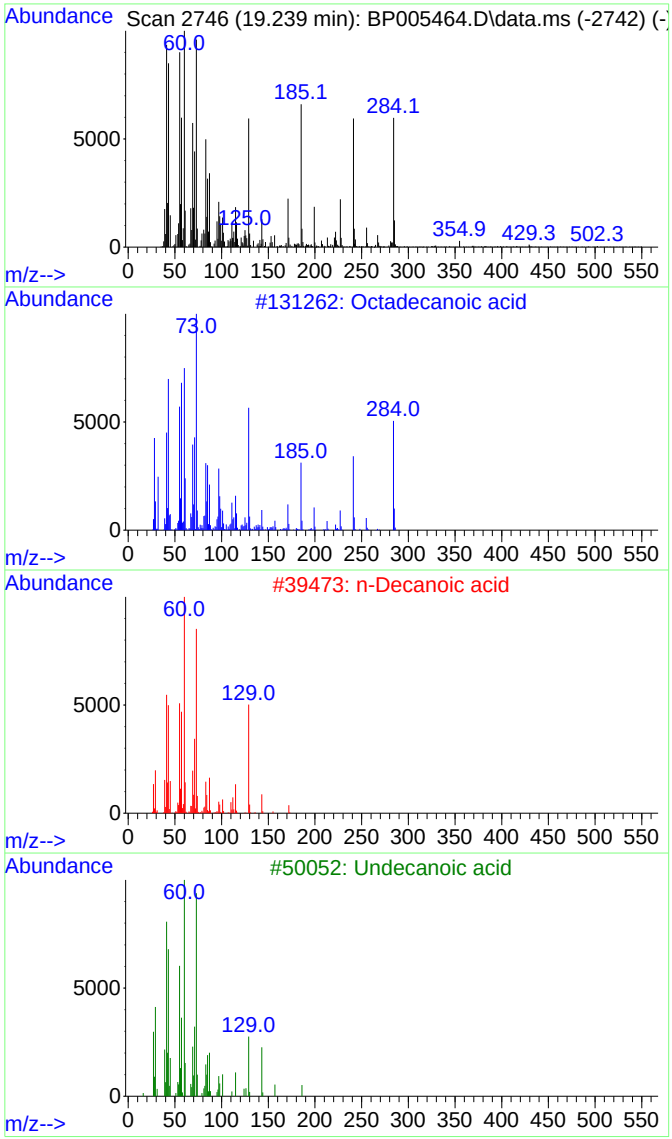
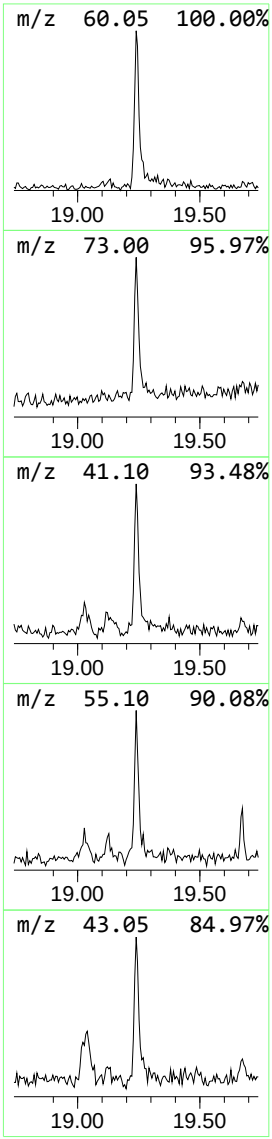
Peak Number 10 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	6.82 ng	85734	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2	n-Decanoic acid	172	C10H20O2	000334-48-5	56
3	Undecanoic acid	186	C11H22O2	000112-37-8	55
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	50
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	46





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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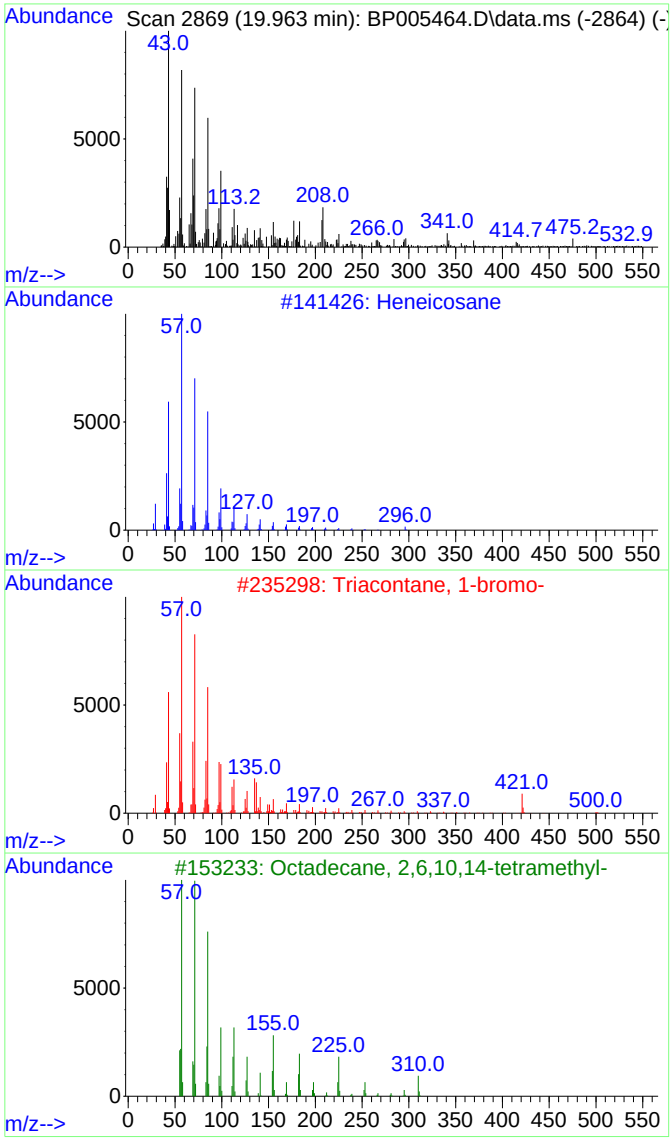
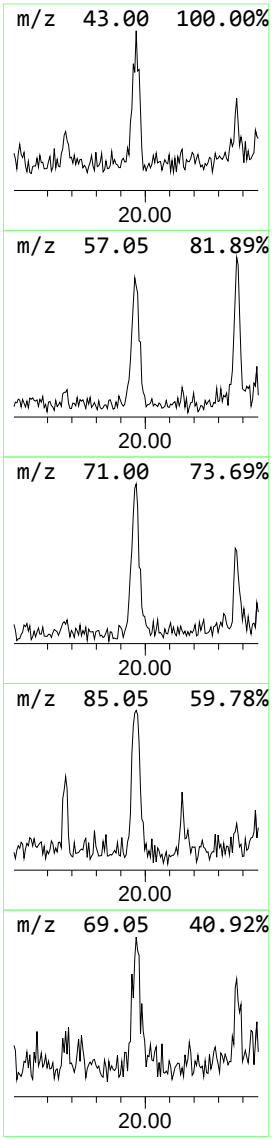
Peak Number 11 Heneicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.963	2.96 ng	37294	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	76
3	Octadecane, 2,6,10,14-tetramethyl-	310	C22H46	054964-82-8	76
4	Tetratetracontane	619	C44H90	007098-22-8	74
5	Dodecane, 4-methyl-	184	C13H28	006117-97-1	68





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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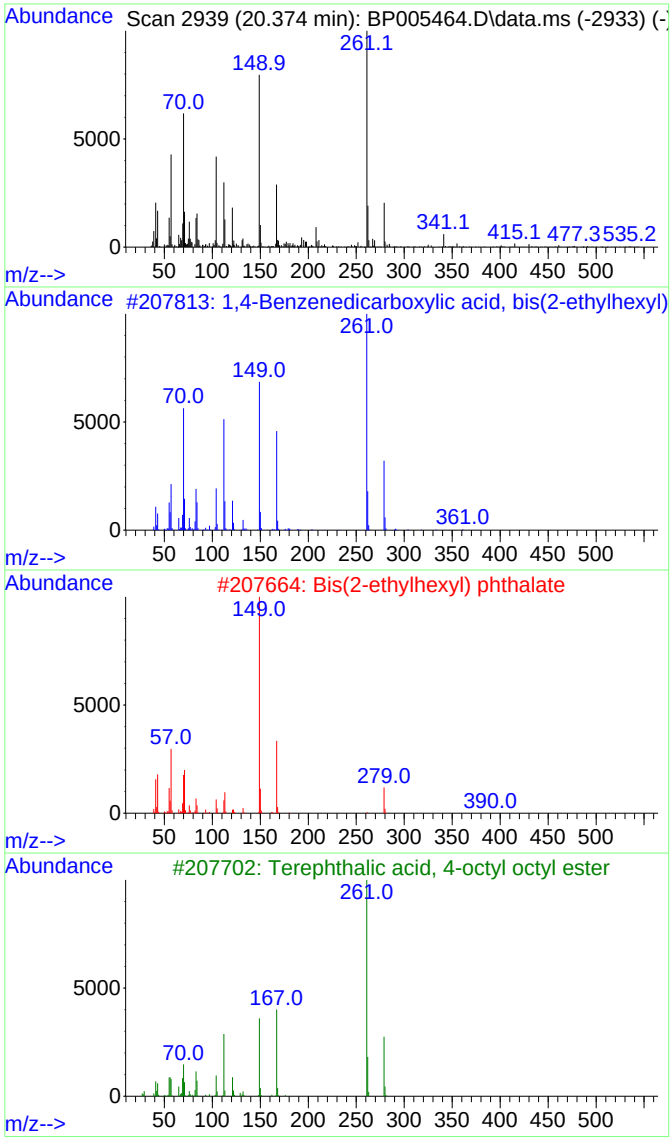
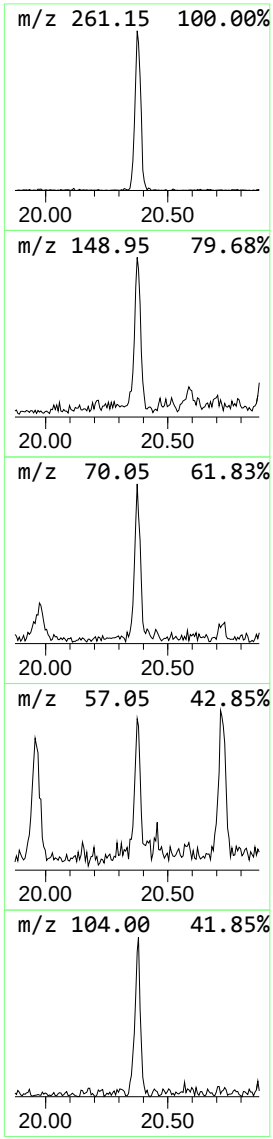
Peak Number 12 1,4-Benzenedicarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.374	5.54 ng	69699	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	58
2	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	42
3	Terephthalic acid, 4-octyl octyl...	390	C24H38O4	1000323-73-7	27
4	Diisooctyl phthalate	390	C24H38O4	000131-20-4	27
5	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	27





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Acq On : 10 May 2021 12:39
Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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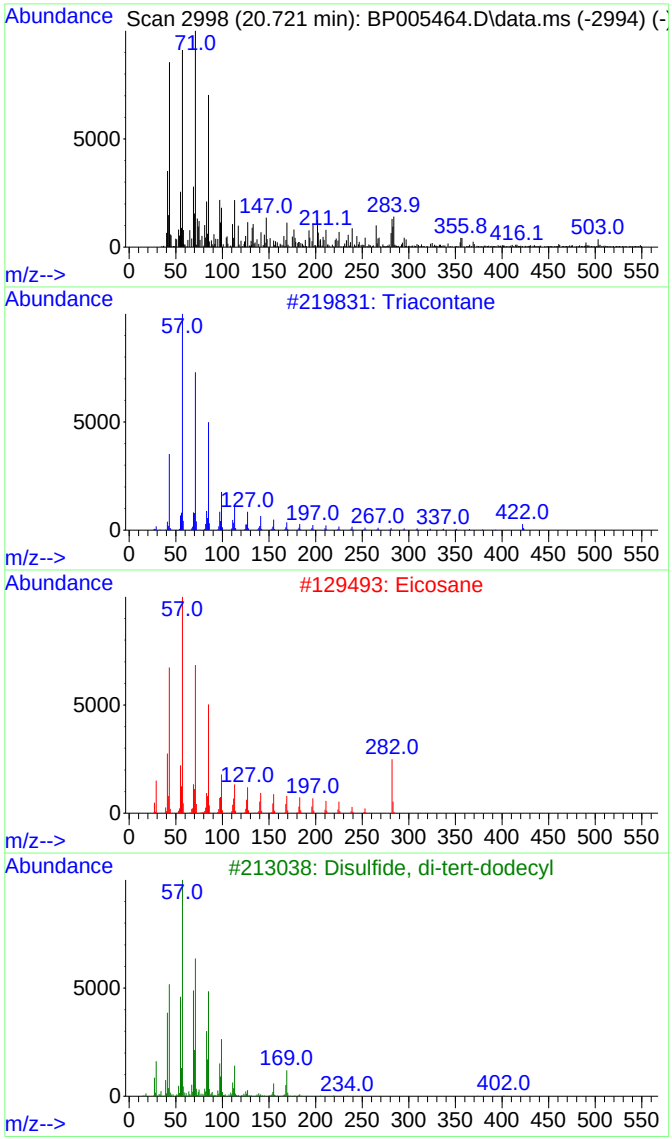
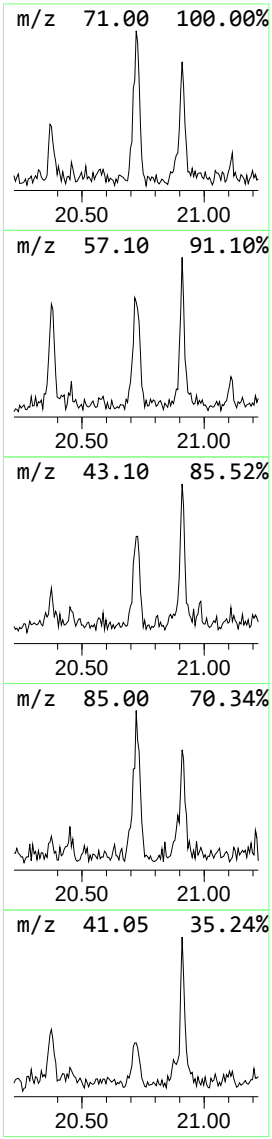
Peak Number 13 Triacontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.721	3.05 ng	38366	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	89
2	Eicosane	282	C20H42	000112-95-8	76
3	Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	74
4	Tetratriacontane	479	C34H70	014167-59-0	72
5	Docosane	310	C22H46	000629-97-0	64





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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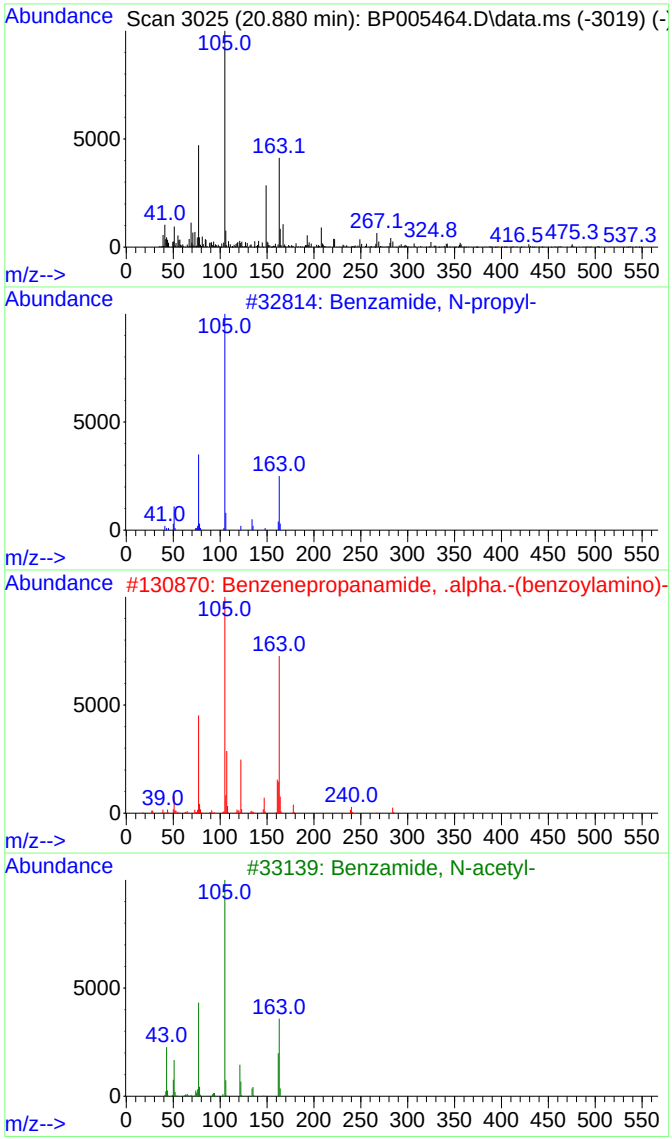
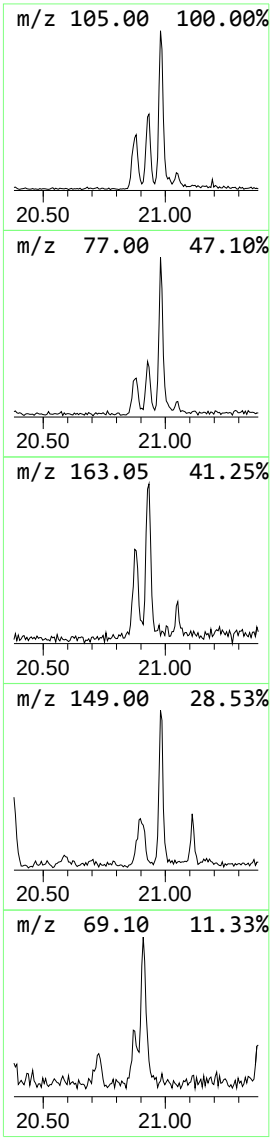
Peak Number 14 unknown20.88 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.880	3.82 ng	48082	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, N-propyl-	163	C10H13NO	010546-70-0	47
2	Benzenepropanamide, .alpha.-(ben...	284	C16H16N2O3	058690-81-6	47
3	Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	47
4	2,3-Dimethyl-4a,8a-diphenyl-hexa...	326	C20H22O4	1000297-07-0	38
5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	35





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

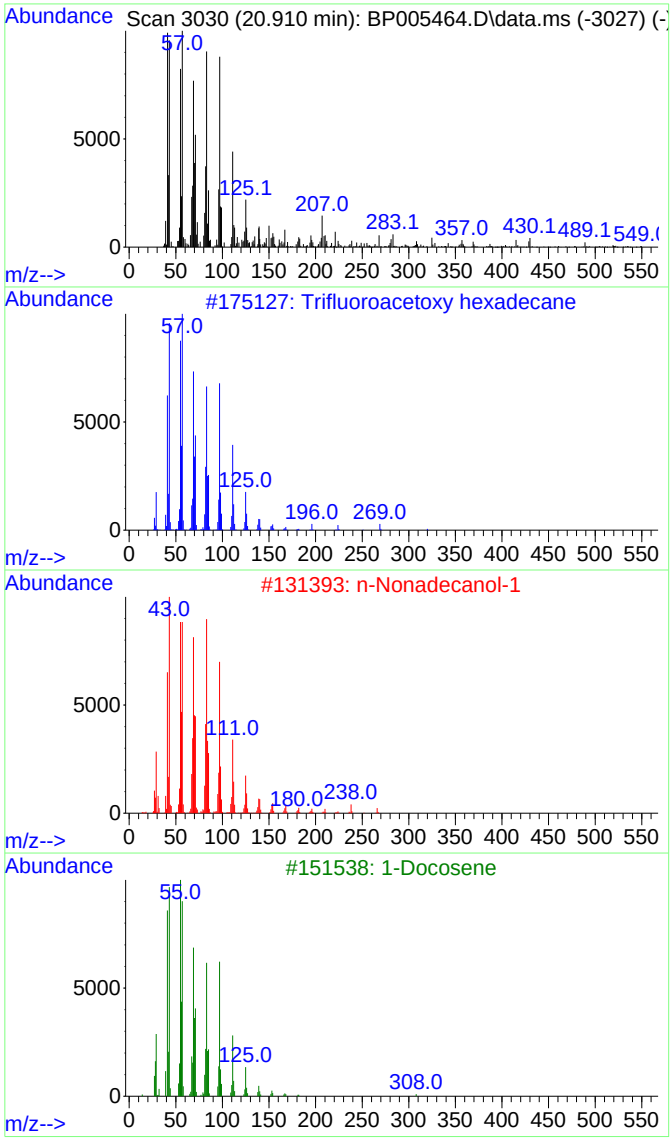
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

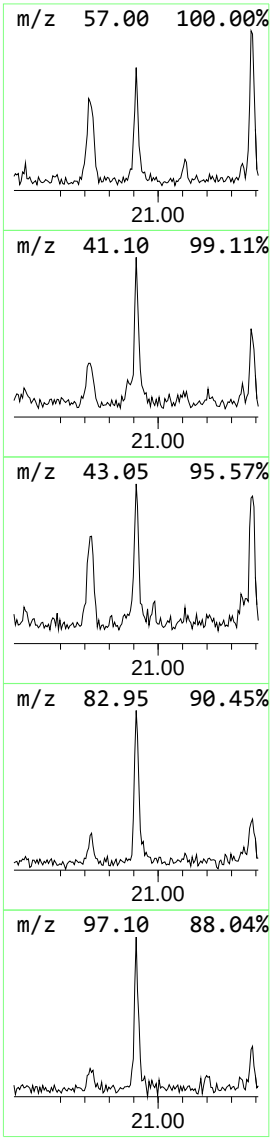
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Trifluoroacetoxy hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.910	8.30 ng	104419	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	90
3	1-Docosene	308	C22H44	001599-67-3	89
4	1-Eicosene	280	C20H40	003452-07-1	89
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	87





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Acq On : 10 May 2021 12:39
Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

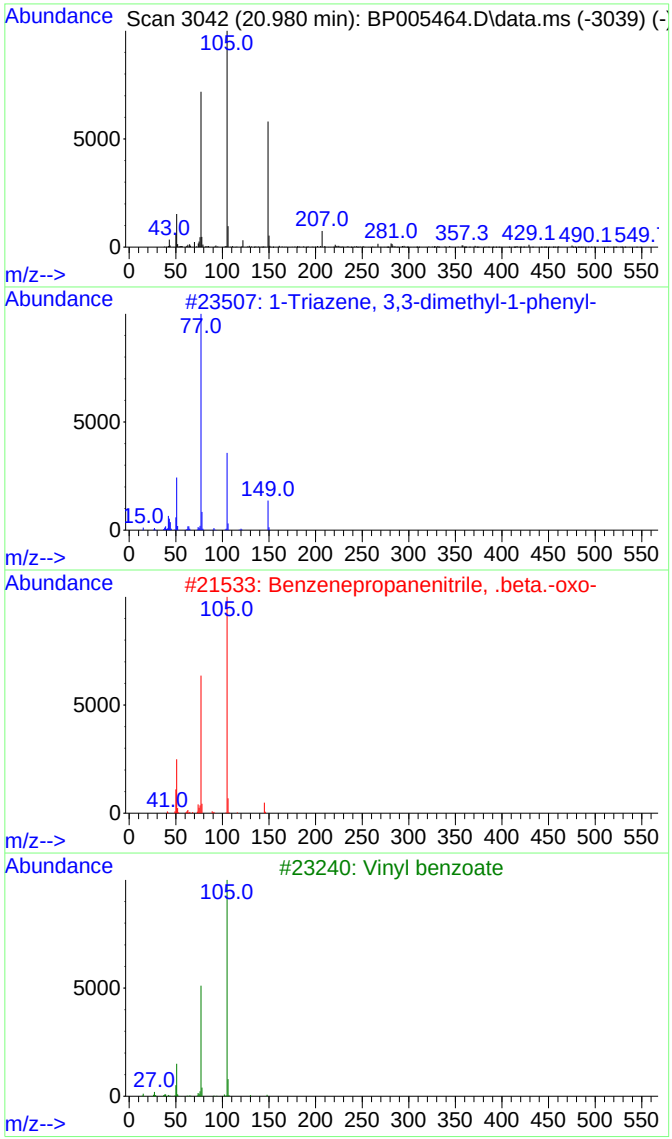
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

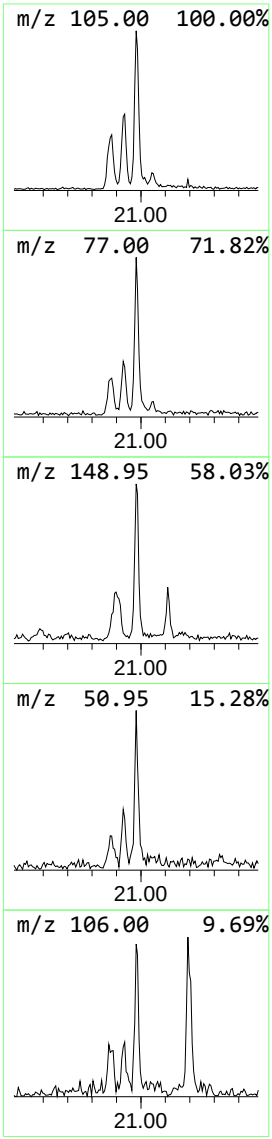
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1-Triazene, 3,3-dimethyl-1-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.980	4.86 ng	61121	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Triazene, 3,3-dimethyl-1-phenyl-	149	C8H11N3	007227-91-0	72
2	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	43
3	Vinyl benzoate	148	C9H8O2	000769-78-8	43
4	Benzoyl bromide	184	C7H5BrO	000618-32-6	43
5	Pentanediamide, N,N'-di-benzoyloxy-	370	C19H18N2O6	1000253-26-3	43





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Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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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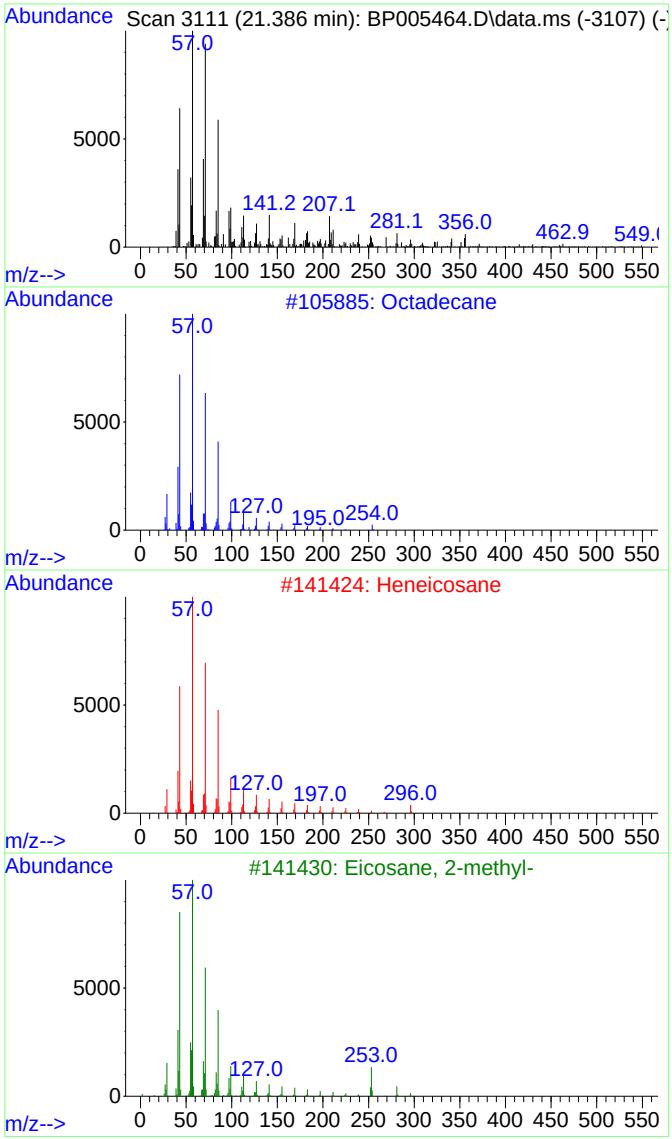
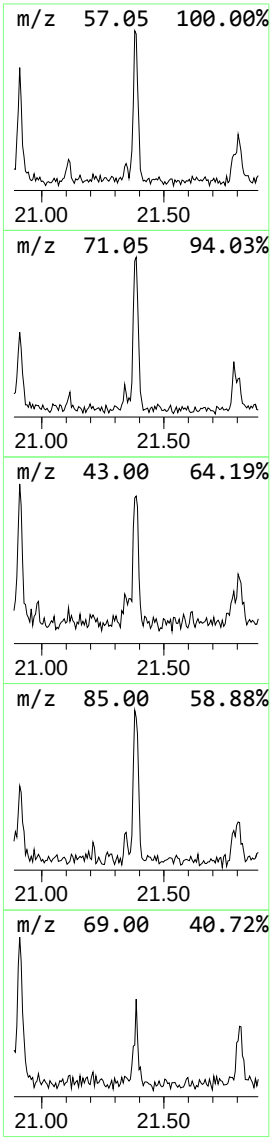
Peak Number 17 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.386	4.04 ng	50871	Chrysene-d12	21.198

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	95
2	Heneicosane	296	C21H44	000629-94-7	91
3	Eicosane, 2-methyl-	296	C21H44	001560-84-5	91
4	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
5	Hexatriacontane	507	C36H74	000630-06-8	91





Tentatively Identified Compound (LSC) summary

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Data File : BP005464.D
Acq On : 10 May 2021 12:39
Operator : CG/JU
Sample : M2306-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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
unknown4.82	4.828	3.2	ng	13303	1	7.705	82190	20.0
4-((1E)-3-Hydro...	16.516	5.8	ng	47054	4	17.104	161869	20.0
Benzamide, N-ac...	17.775	8.0	ng	64850	4	17.104	161869	20.0
Ethanone, 2-hyd...	17.927	9.8	ng	79034	4	17.104	161869	20.0
n-Decanoic acid	18.004	16.8	ng	136181	4	17.104	161869	20.0
3,5-Dimethoxy-4...	18.280	9.8	ng	78871	4	17.104	161869	20.0
2-Pentanone, 1-...	18.328	25.6	ng	207637	4	17.104	161869	20.0
Octacosane	19.027	3.0	ng	24627	4	17.104	161869	20.0
3(2H)-Pyridazin...	19.127	2.1	ng	16770	4	17.104	161869	20.0
Octadecanoic acid	19.239	6.8	ng	85734	5	21.198	251568	20.0
Heneicosane	19.963	3.0	ng	37294	5	21.198	251568	20.0
1,4-Benzenedica...	20.374	5.5	ng	69699	5	21.198	251568	20.0
Triacontane	20.721	3.0	ng	38366	5	21.198	251568	20.0
unknown20.88	20.880	3.8	ng	48082	5	21.198	251568	20.0
Trifluoroacetox...	20.910	8.3	ng	104419	5	21.198	251568	20.0
1-Triazene, 3,3...	20.980	4.9	ng	61121	5	21.198	251568	20.0
Octadecane	21.386	4.0	ng	50871	5	21.198	251568	20.0