

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
 Data File : BG048707.D
 Acq On : 15 Apr 2021 13:40
 Operator : CG/JU
 Sample : M2015-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CL-02-041221

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG048707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.635	384	391	403	rBV	392675	626444	9.83%	3.112%
2	7.244	658	665	677	rVB	404065	698504	10.96%	3.470%
3	8.085	800	808	816	rBV	101058	164084	2.58%	0.815%
4	9.254	998	1007	1014	rBV	257014	458458	7.20%	2.277%
5	10.911	1282	1289	1296	rVB	130518	225832	3.54%	1.122%
6	13.355	1698	1705	1712	rBV	607945	948643	14.89%	4.712%
7	13.619	1744	1750	1758	rBV2	60278	102038	1.60%	0.507%
8	14.742	1934	1941	1946	rBV2	187604	295219	4.63%	1.466%
9	15.846	2122	2129	2135	rBV2	38341	66353	1.04%	0.330%
10	16.228	2187	2194	2203	rBV	525744	827106	12.98%	4.108%
11	17.491	2402	2409	2415	rBV	231299	373555	5.86%	1.856%
12	18.373	2554	2559	2564	rBV5	75121	108722	1.71%	0.540%
13	19.072	2666	2678	2688	rBV2	632385	2076648	32.59%	10.315%
14	19.213	2688	2702	2710	rBV	843376	2529884	39.71%	12.567%
15	19.324	2710	2721	2730	rVB	2331593	6371514	100.00%	31.649%
16	19.465	2737	2745	2753	rVB	188704	409636	6.43%	2.035%
17	19.642	2769	2775	2781	rVB2	57479	108880	1.71%	0.541%
18	20.106	2849	2854	2858	rBV	1015840	1291063	20.26%	6.413%
19	20.153	2858	2862	2867	rVB4	60352	102674	1.61%	0.510%
20	20.905	2984	2990	2997	rVB2	99657	186821	2.93%	0.928%
21	21.416	3072	3077	3087	rBV3	63779	136740	2.15%	0.679%
22	21.704	3121	3126	3132	rVB2	94789	175573	2.76%	0.872%
23	21.792	3135	3141	3147	rVV	214802	374198	5.87%	1.859%
24	22.621	3275	3282	3292	rVB3	105514	253155	3.97%	1.257%
25	23.666	3453	3460	3470	rVB3	88084	242757	3.81%	1.206%
26	24.900	3660	3670	3679	rBV5	77865	235101	3.69%	1.168%
27	25.135	3700	3710	3719	rBV2	129470	405388	6.36%	2.014%
28	26.351	3907	3917	3927	rBV5	54164	191385	3.00%	0.951%
29	28.132	4213	4220	4231	rVB7	39081	145348	2.28%	0.722%

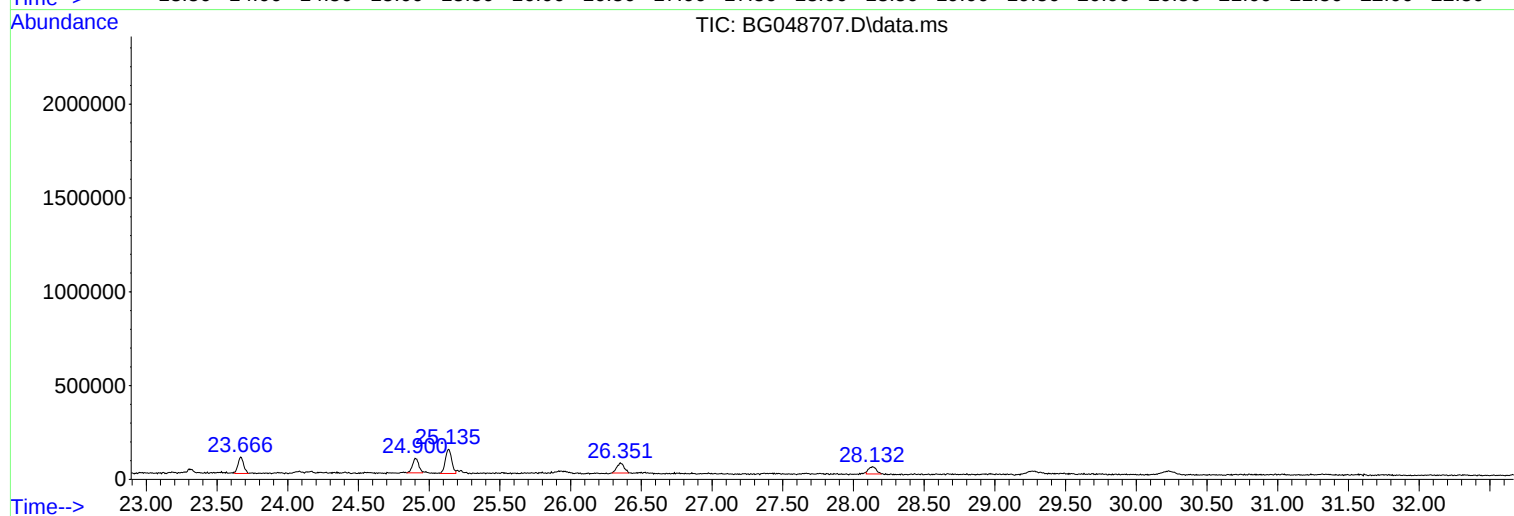
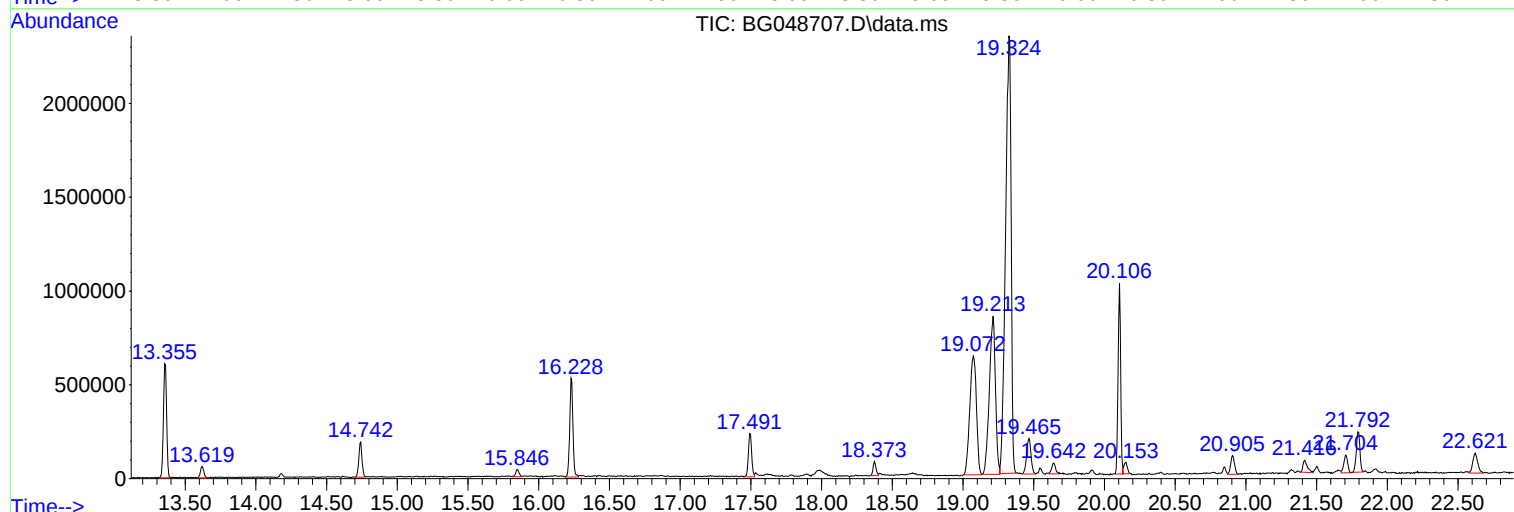
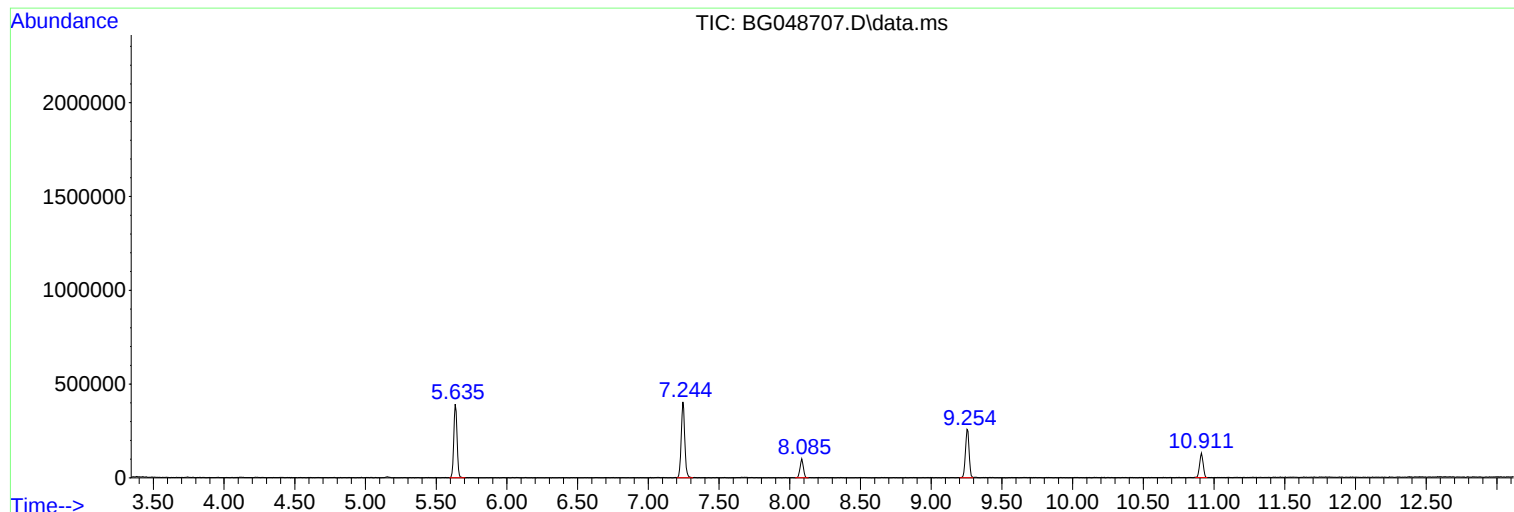
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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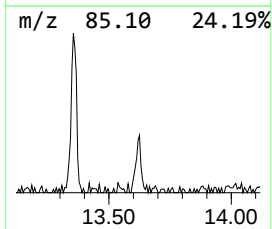
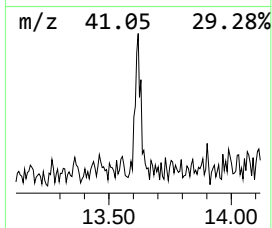
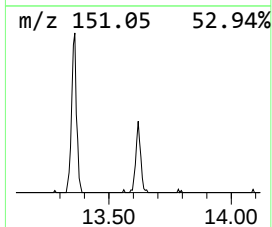
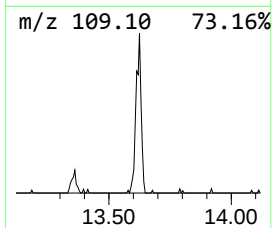
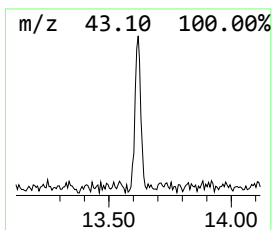
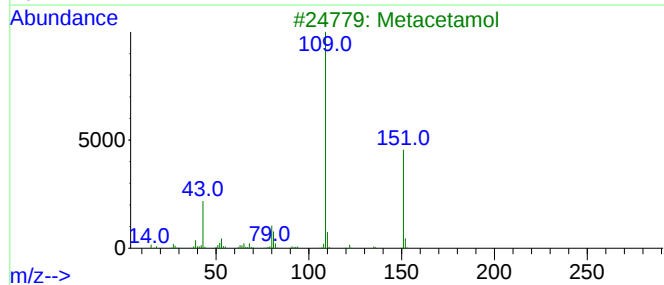
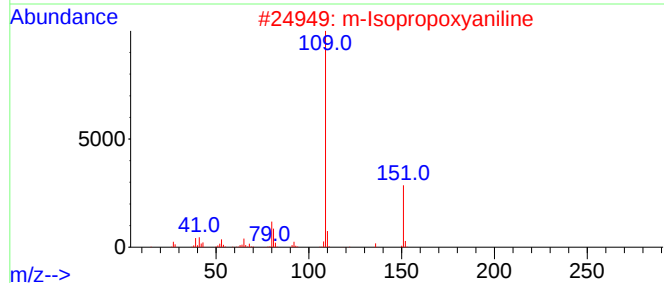
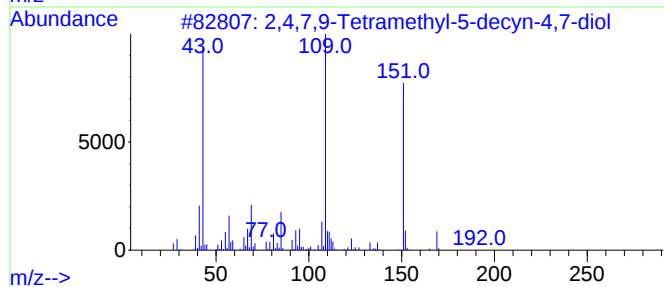
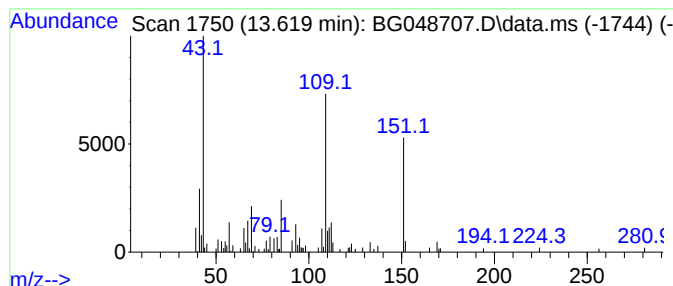
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.619	6.91 ng	102038	Acenaphthene-d10	14.736

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	72		
2	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	58		
3	Metacetamol	151	C8H9NO2	000621-42-1	53		
4	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53		
5	(-)-10-Camphorsulfonyl chloride	250	C10H15ClO3S	039262-22-1	47		



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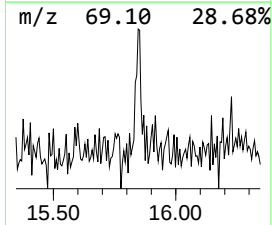
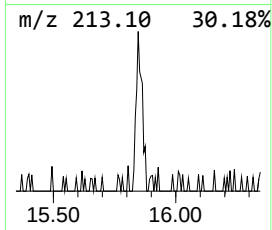
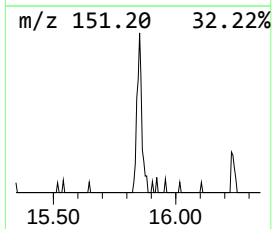
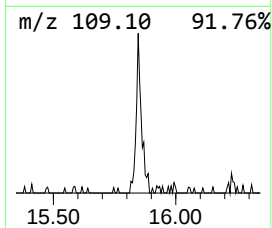
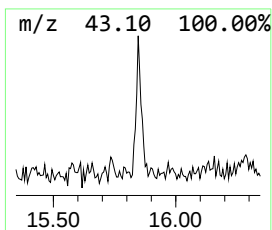
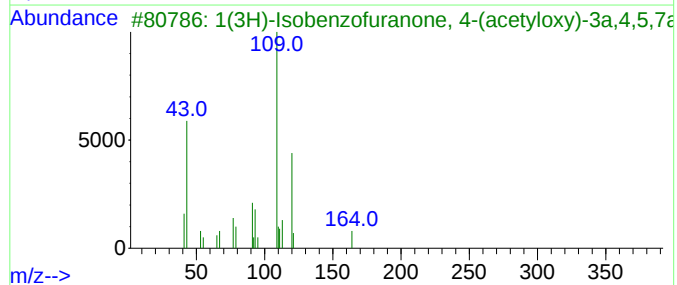
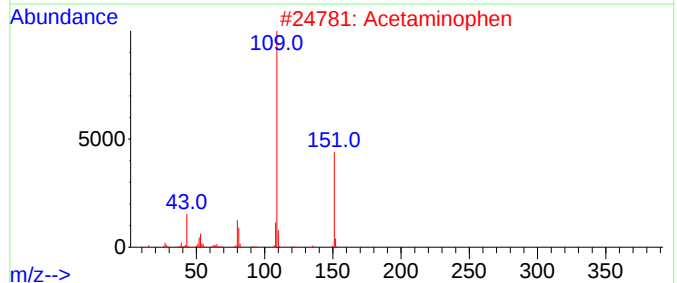
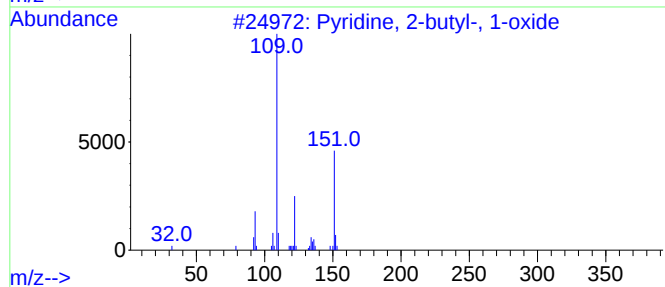
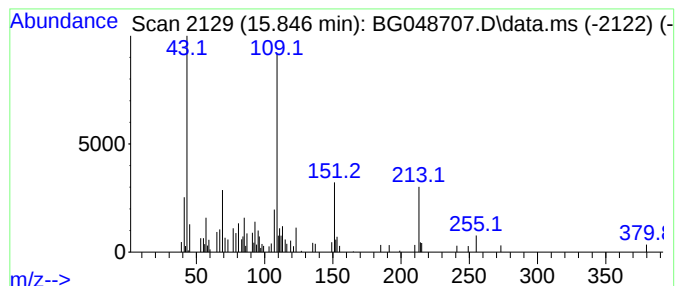
Instrument :
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ClientSampleId :
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Peak Number 2 unknown15.846 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.846	4.50 ng	66353	Acenaphthene-d10	14.736	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	43
2	Acetaminophen	151	C8H9NO2	000103-90-2	43
3	1(3H)-Isobenzofuranone, 4-(acety...	224	C12H16O4	054346-07-5	43
4	Metacetamol	151	C8H9NO2	000621-42-1	38
5	Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	38



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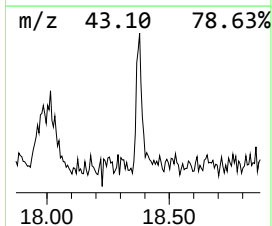
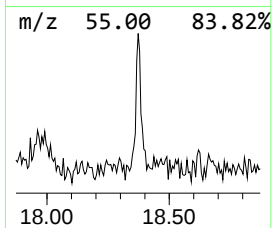
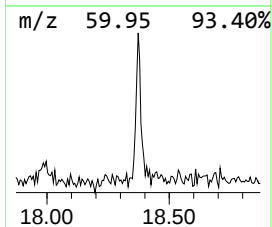
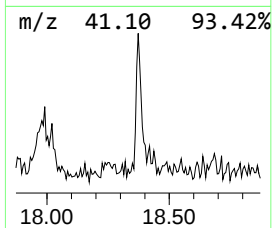
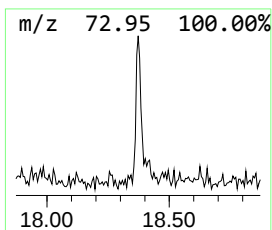
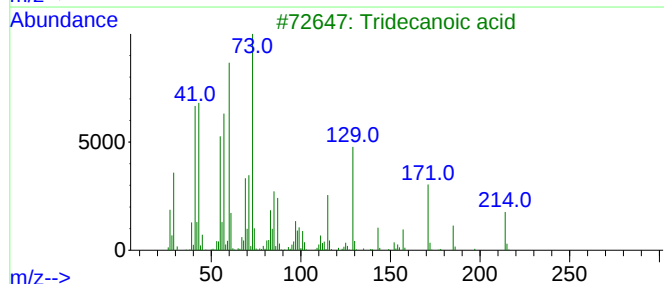
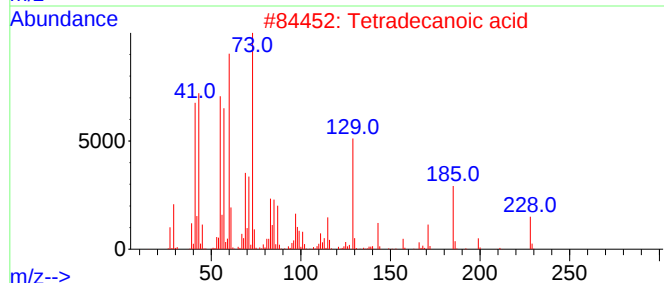
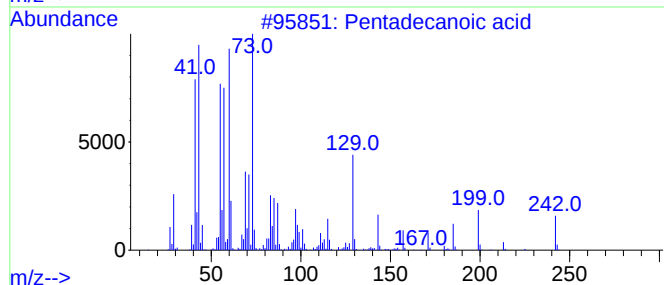
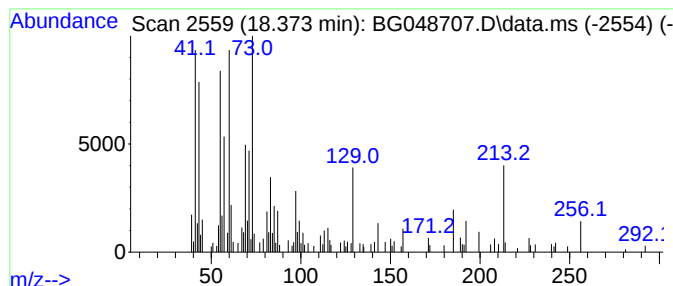
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Pentadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.372	5.82 ng	108722	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid	242	C15H30O2	001002-84-2	95
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	70
3			Tridecanoic acid	214	C13H26O2	000638-53-9	70
4			n-Decanoic acid	172	C10H20O2	000334-48-5	58
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55



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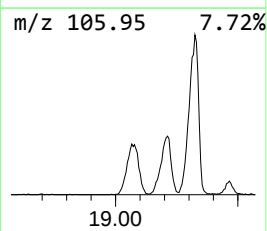
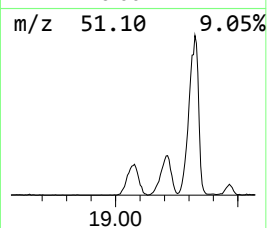
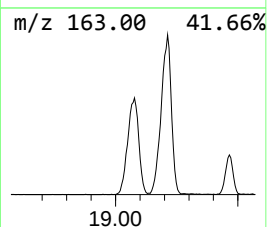
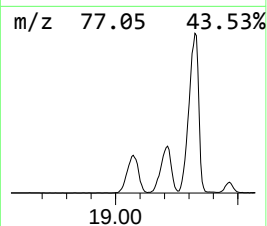
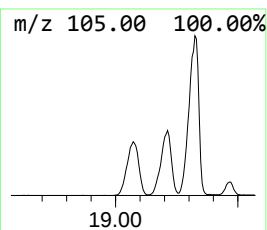
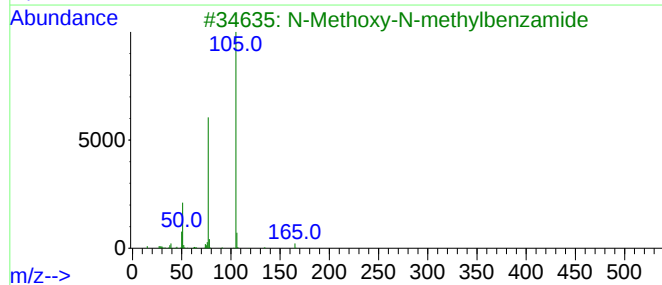
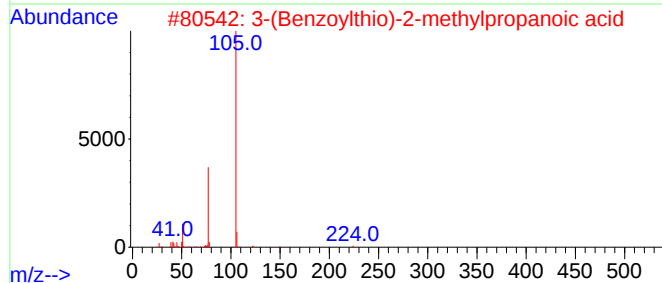
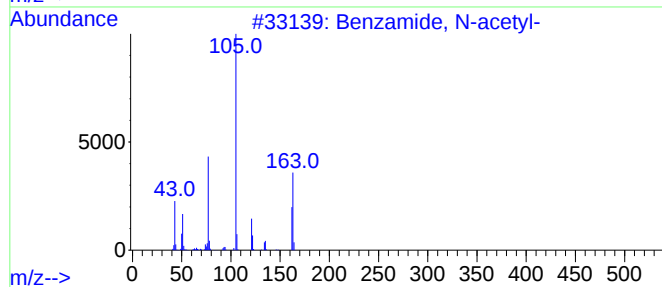
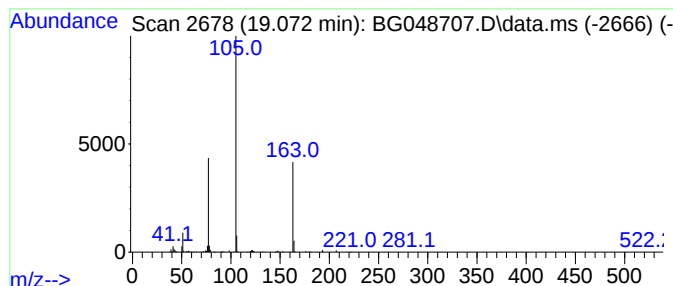
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Peak Number 4 Benzamide, N-acetyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.072	111.18 ng	2076650	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	56
2			3-(Benzoylthio)-2-methylpropanoi...	224	C11H12O3S	067714-34-5	47
3			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47
4			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	47
5			Benzoic acid, pentafluorophenyl ...	288	C13H5F5O2	1000360-67-9	47



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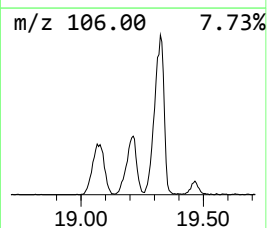
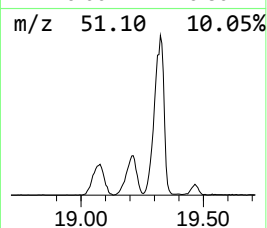
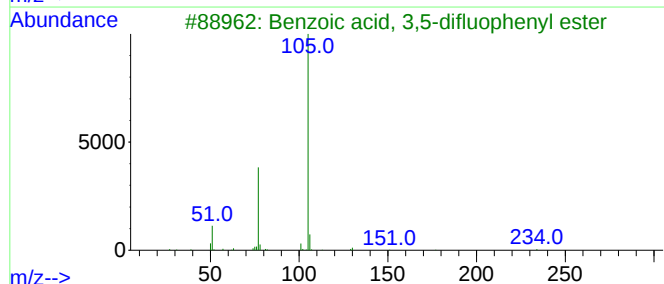
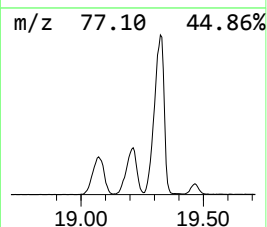
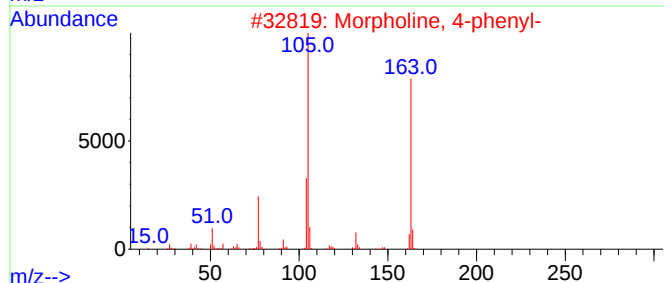
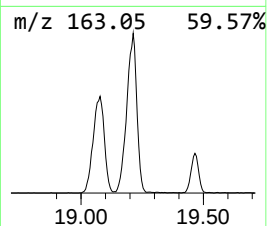
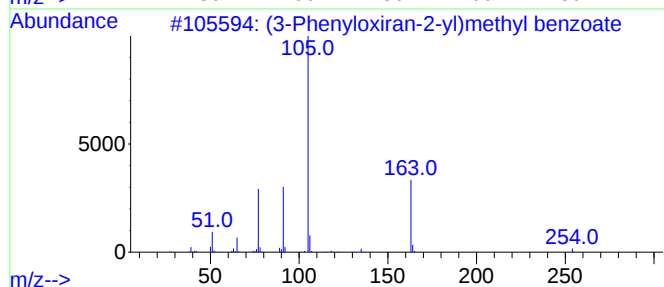
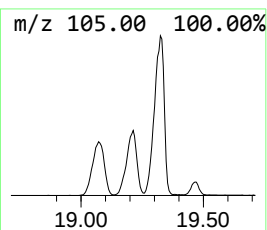
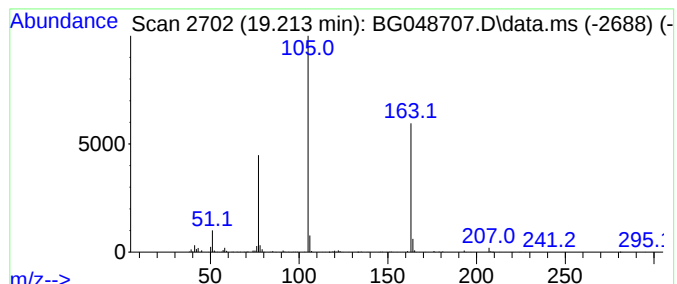
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Peak Number 5 (3-Phenyloxiran-2-yl)methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.213	135.45 ng	2529880	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	59
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
3			Benzoic acid, 3,5-difluophenyl e...	234	C13H8F2O2	1000357-79-5	47
4			Benzo[b]thiophene-3-carboxylic a...	343	C18H17NO4S	096334-43-9	47
5			Vinyl benzoate	148	C9H8O2	000769-78-8	47



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CL-02-041221

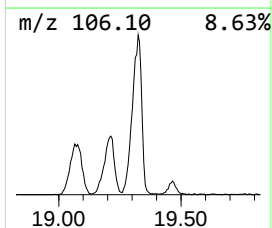
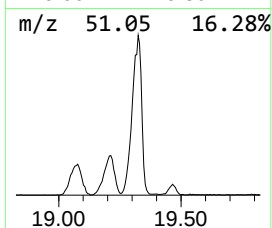
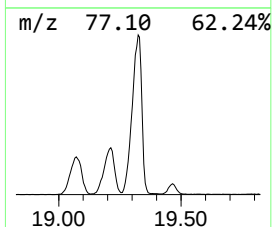
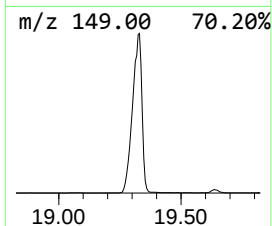
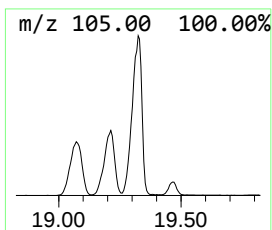
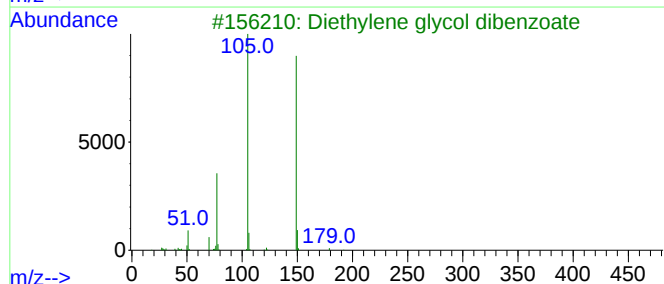
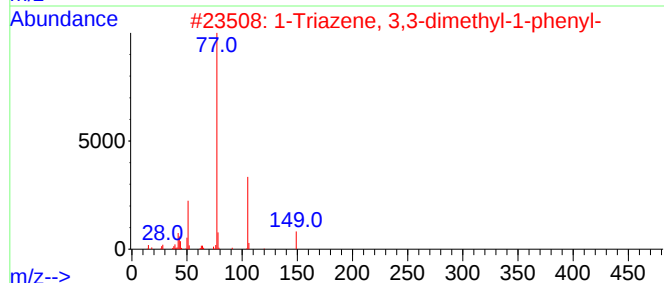
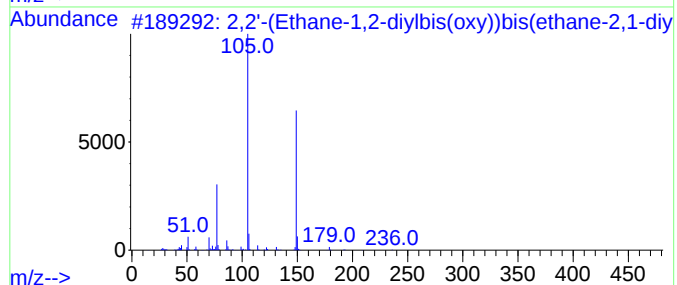
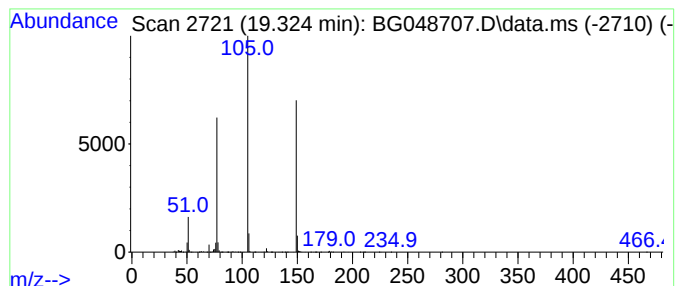
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.324	341.13 ng	6371510	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	74		
2	1-Triazene, 3,3-dimethyl-1-phenyl-	149	C8H11N3	007227-91-0	72		
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	53		
4	Vinyl benzoate	148	C9H8O2	000769-78-8	47		
5	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

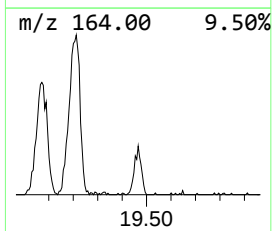
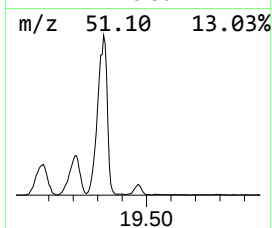
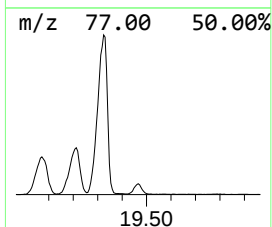
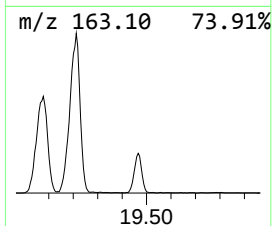
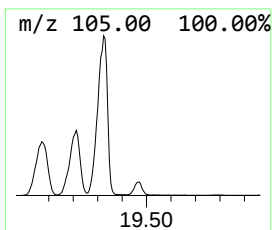
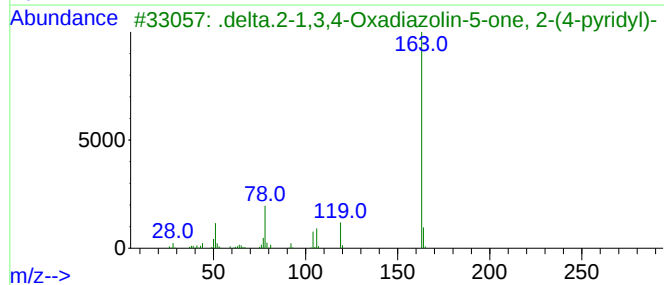
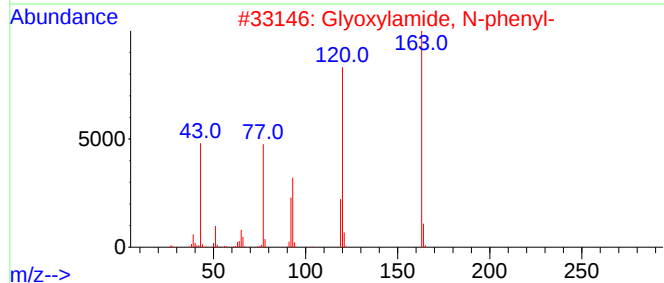
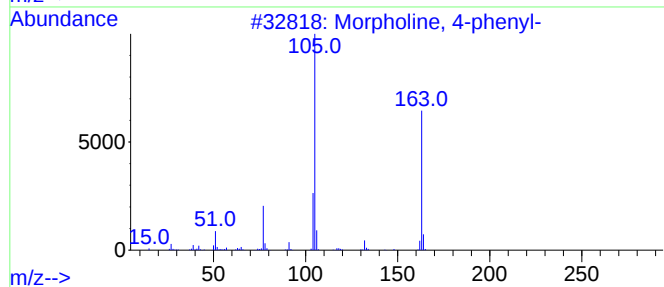
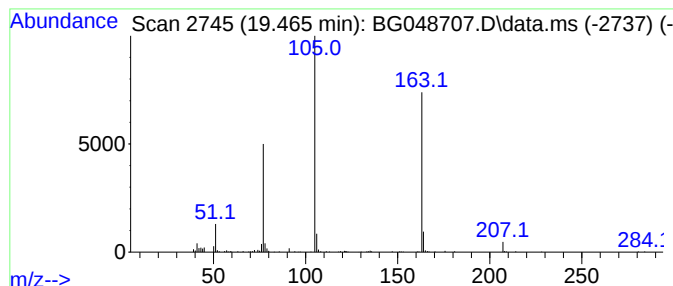
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Morpholine, 4-phenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.465	21.93 ng	409636	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	80
2			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	59
3			.delta.2-1,3,4-Oxadiazolin-5-one...	163	C7H5N3O2	002845-82-1	47
4			Phthalic acid, 3,4-dimethylpheny...	284	C17H16O4	1000315-69-7	43
5			Propanoic acid, 2-(benzoylamino)...	251	C13H17NO4	141627-52-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
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Instrument :
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ClientSampleId :
CL-02-041221

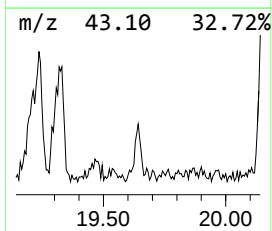
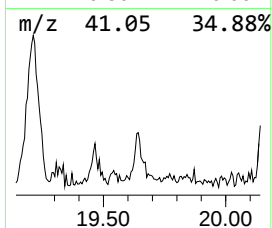
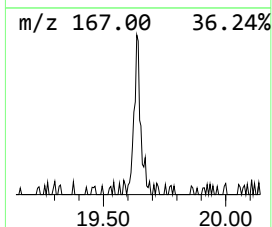
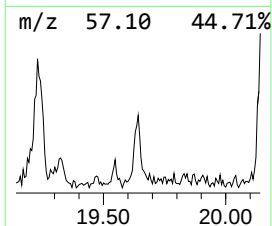
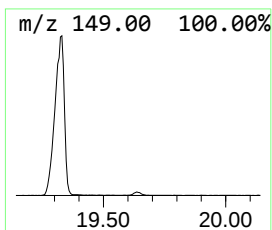
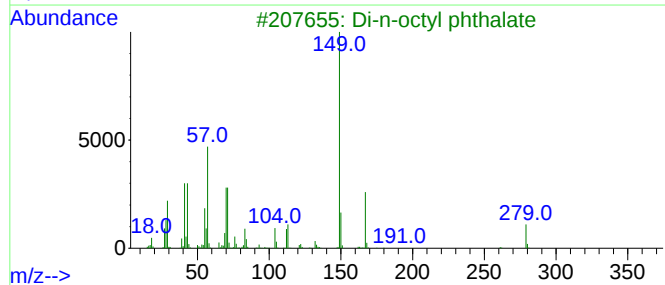
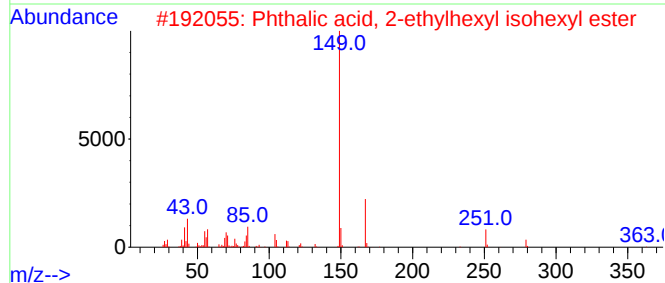
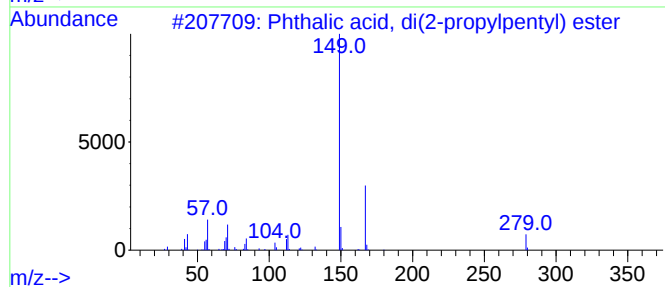
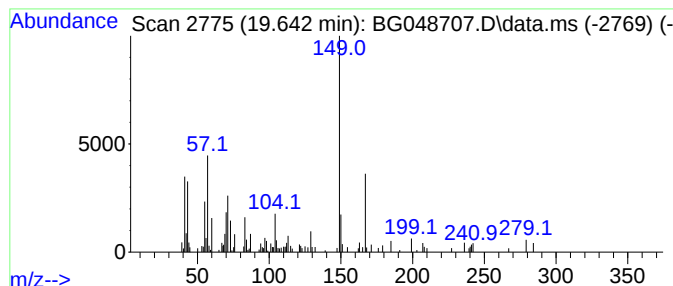
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phthalic acid, di(2-propylp... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.642	5.83 ng	108880	Phenanthrene-d10	17.491

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, di(2-propylpentyl...	390	C24H38O4	1000377-93-5	53
2			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	53
3			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	50
4			2-(Nonyloxycarbonyl)benzoic acid	292	C17H24O4	1000373-89-9	50
5			Phthalic acid, cyclohexyl neopen...	318	C19H26O4	1000315-54-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
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ClientSampleId :
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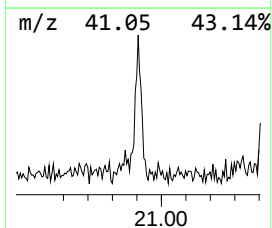
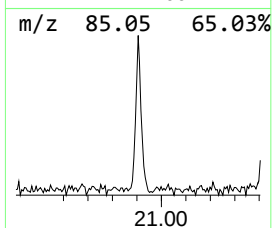
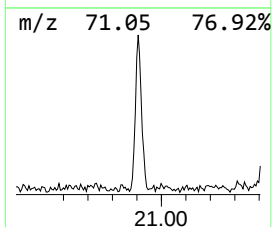
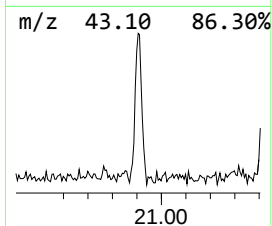
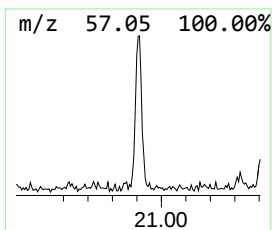
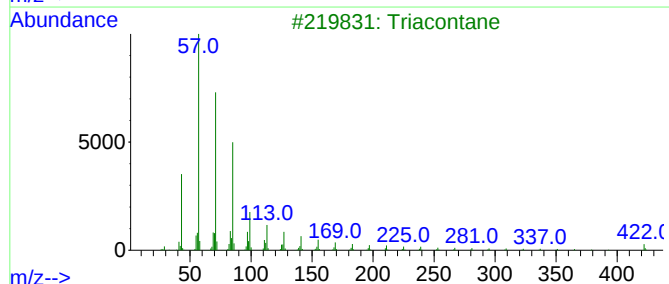
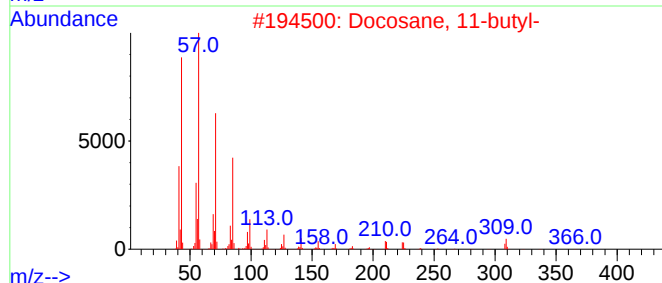
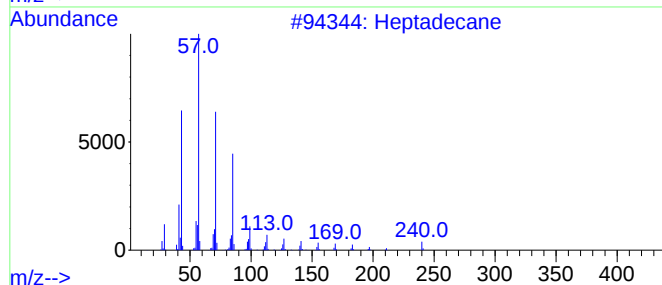
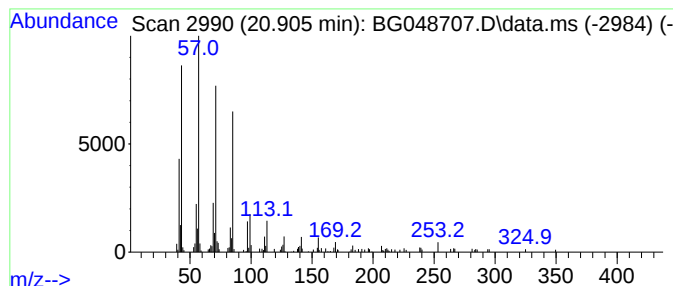
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.905	9.99 ng	186821	Chrysene-d12	21.792

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	93
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Docosane, 11-decyl-	451	C32H66	055401-55-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
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Instrument :
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ClientSampleId :
CL-02-041221

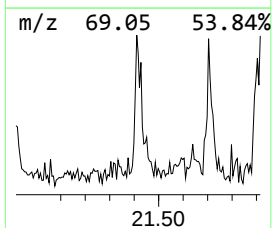
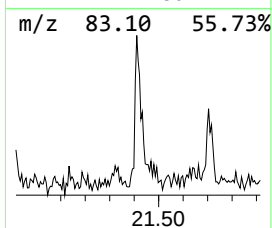
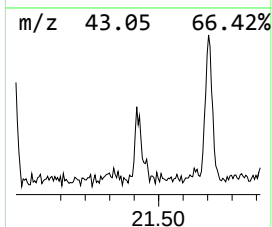
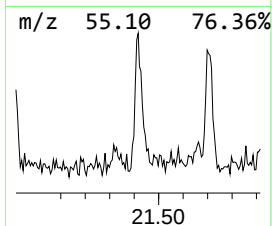
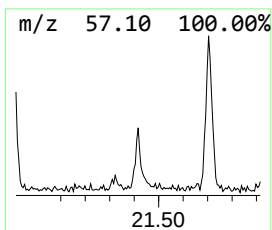
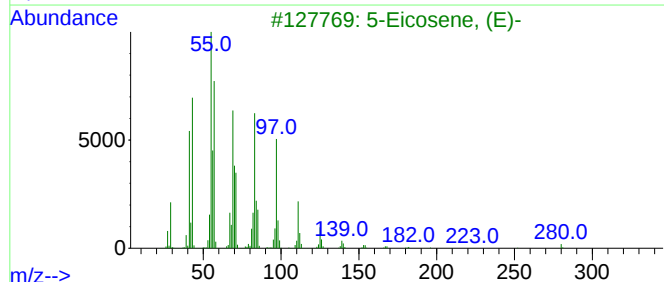
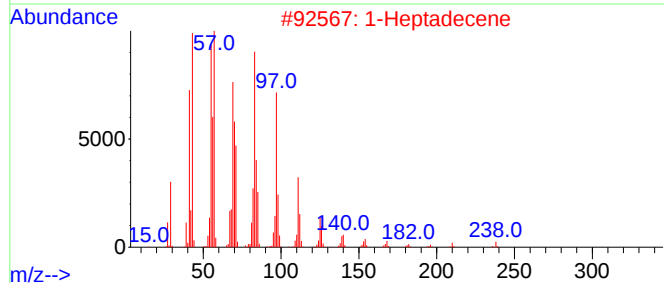
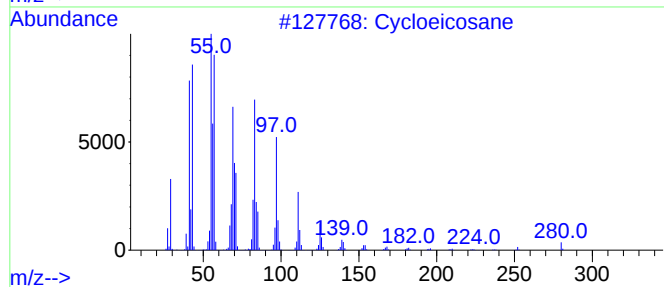
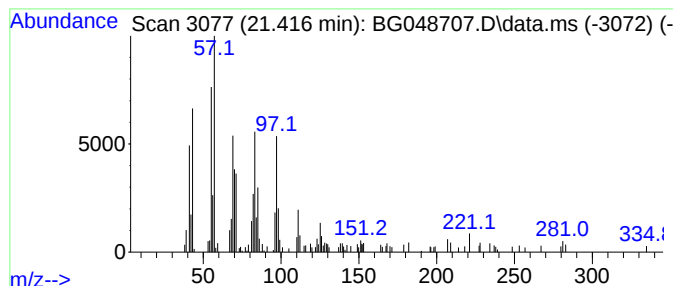
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cycloeicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.416	7.31 ng	136740	Chrysene-d12	21.792

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cycloeicosane	280	C20H40	000296-56-0	92
2		1-Heptadecene	238	C17H34	006765-39-5	91
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	81
4		Cyclotetradecane	196	C14H28	000295-17-0	70
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
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Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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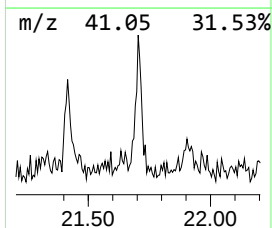
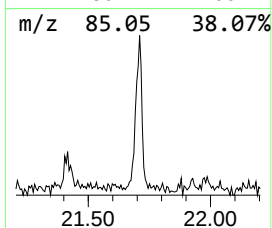
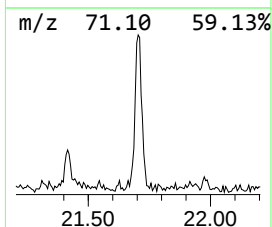
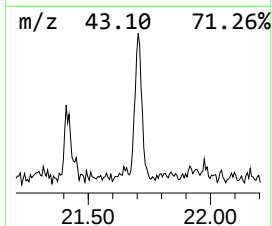
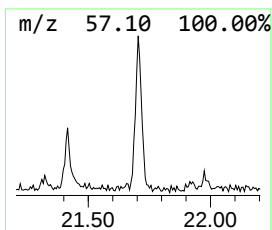
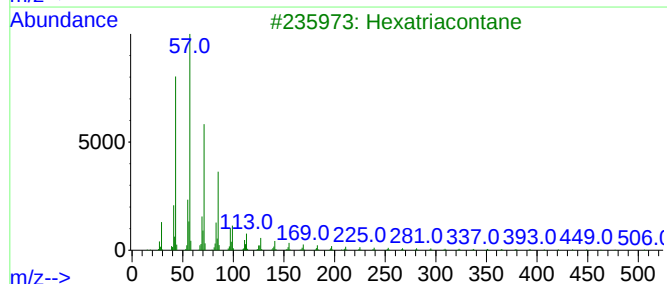
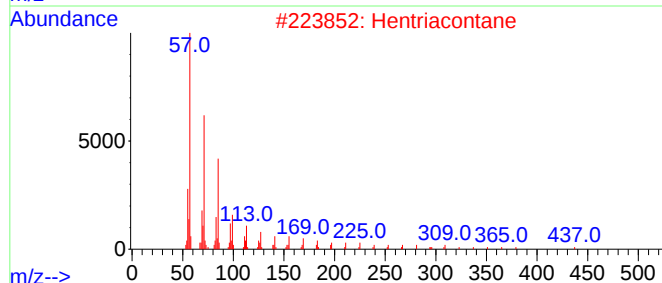
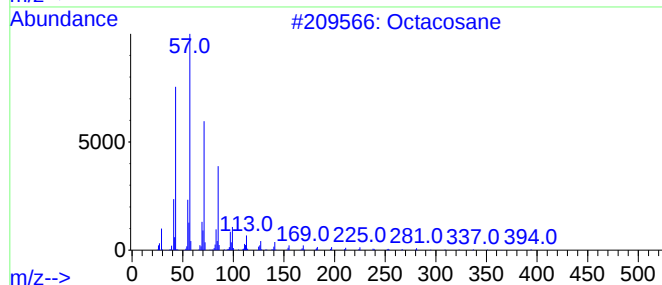
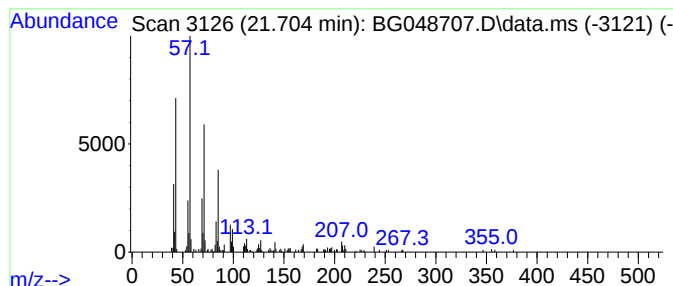
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.704	9.38 ng	175573	Chrysene-d12	21.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Hentriacontane	437	C31H64	000630-04-6	87
3			Hexatriacontane	507	C36H74	000630-06-8	86
4			Nonadecane	268	C19H40	000629-92-5	83
5			Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
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ClientSampleId :
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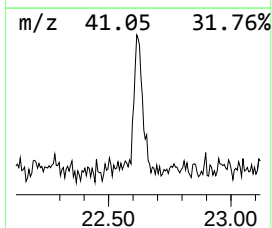
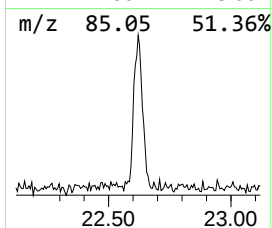
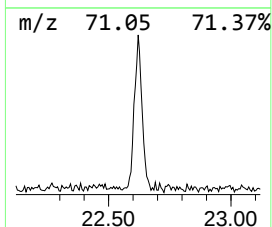
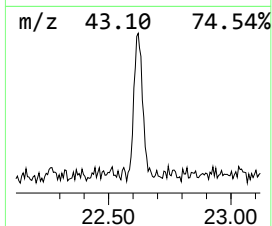
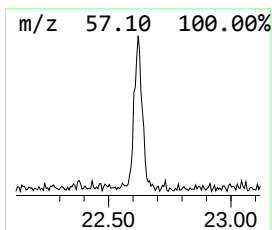
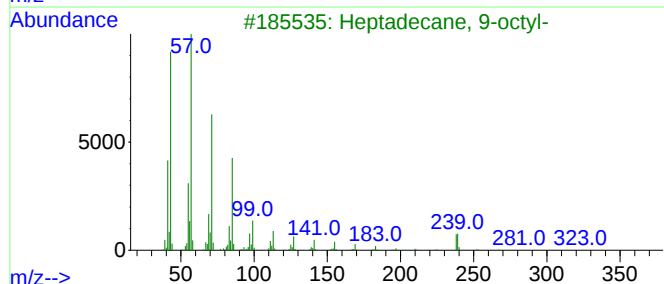
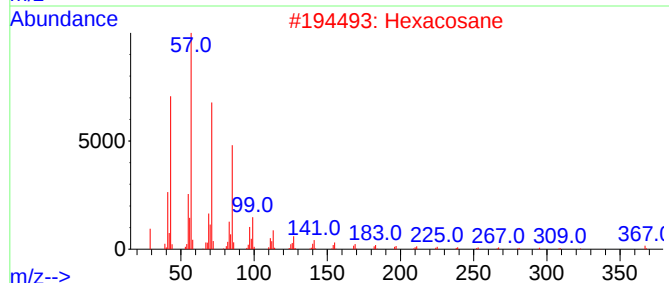
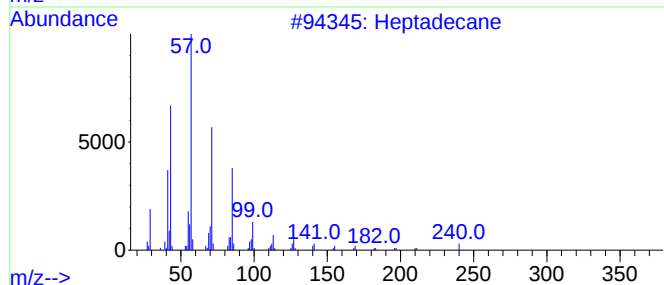
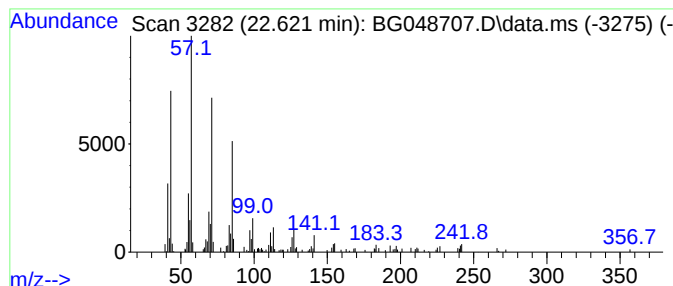
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.621	13.53 ng	253155	Chrysene-d12	21.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	93
2			Hexacosane	366	C26H54	000630-01-3	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

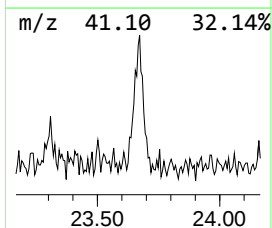
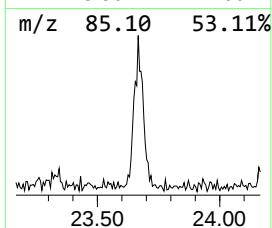
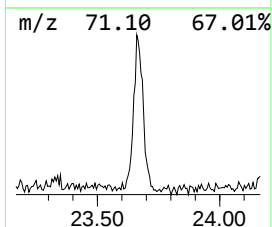
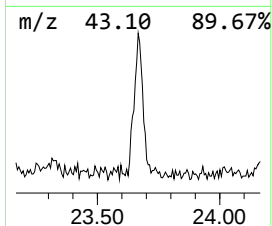
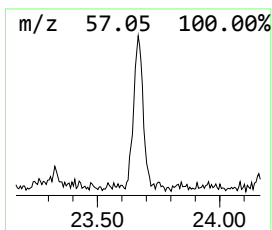
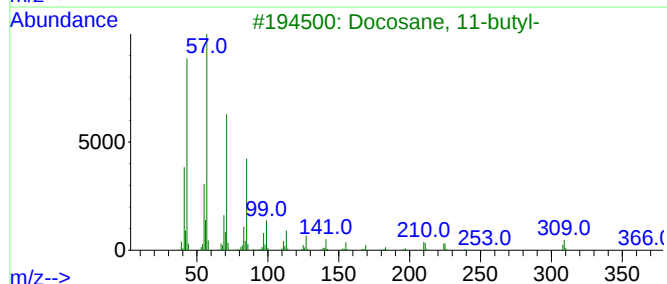
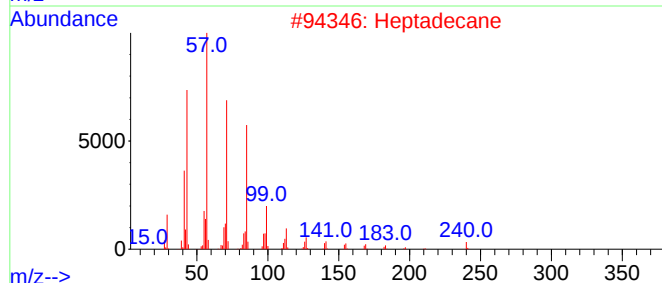
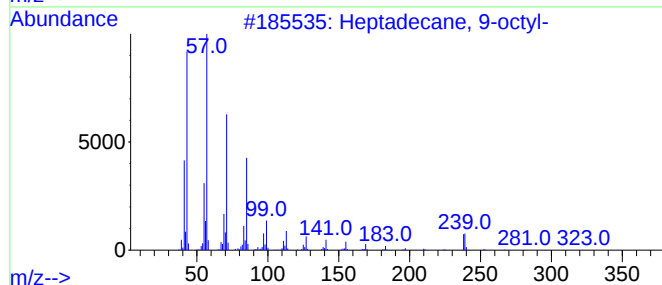
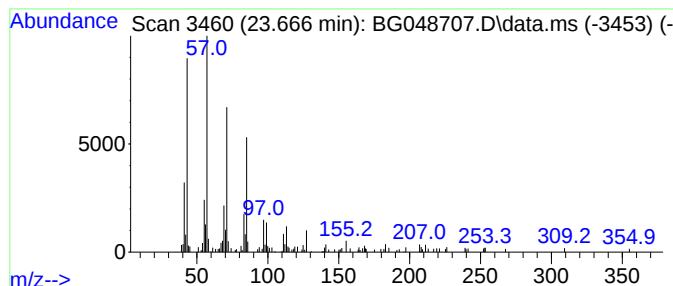
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 9-octyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.666	11.98 ng	242757	Perylene-d12	25.135

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2			Heptadecane	240	C17H36	000629-78-7	91
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4			Docosane, 11-decyl-	451	C32H66	055401-55-3	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

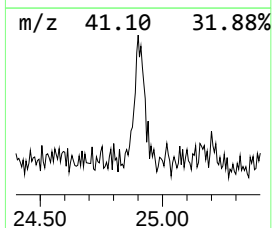
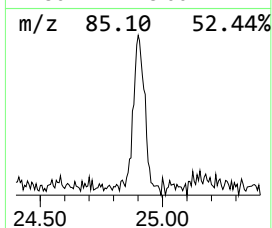
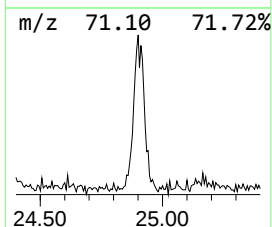
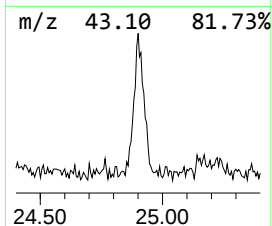
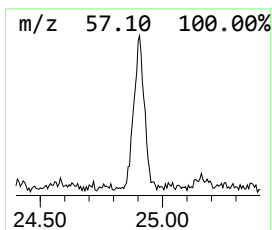
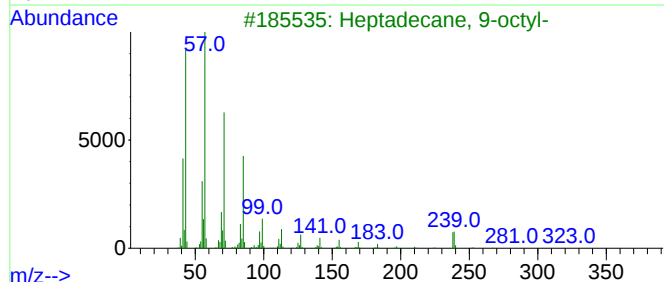
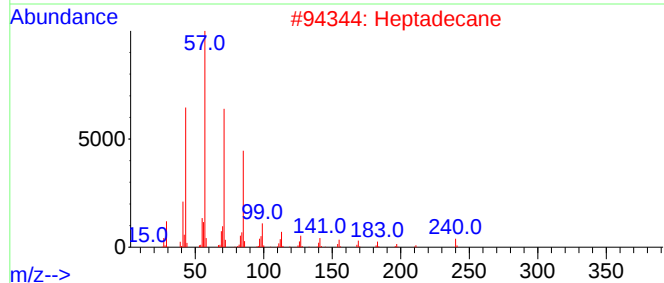
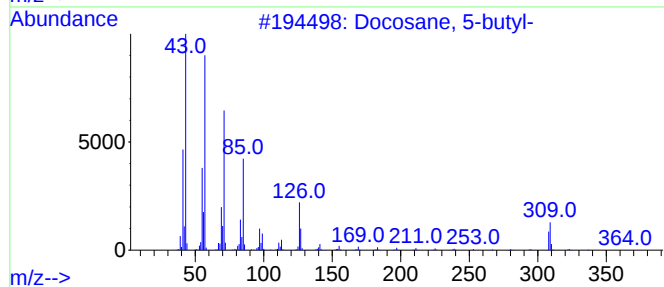
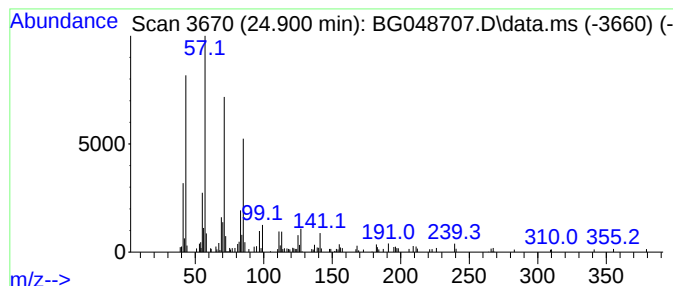
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Docosane, 5-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.900	11.60 ng	235101	Perylene-d12	25.135

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 5-butyl-	366	C26H54	055282-16-1	90
2			Heptadecane	240	C17H36	000629-78-7	89
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4			Tetratetracontane	619	C44H90	007098-22-8	87
5			Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

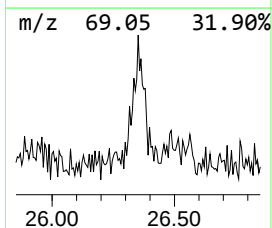
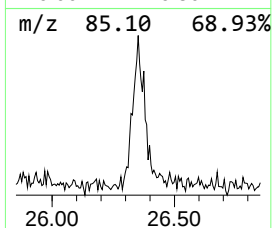
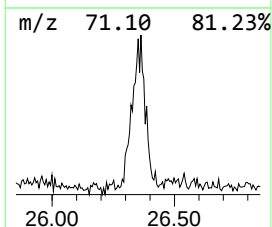
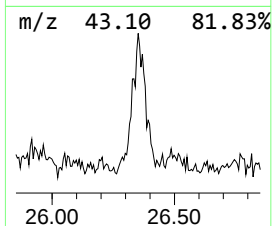
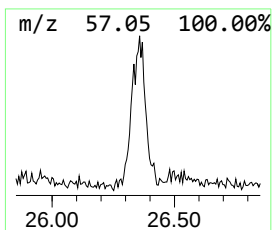
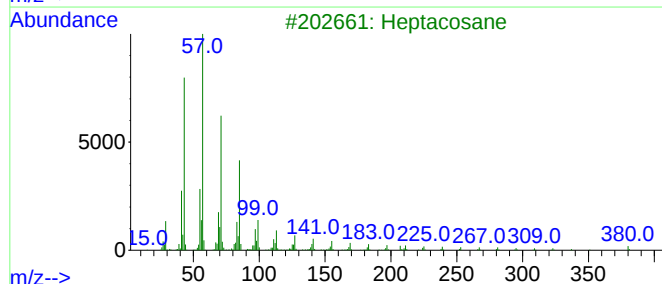
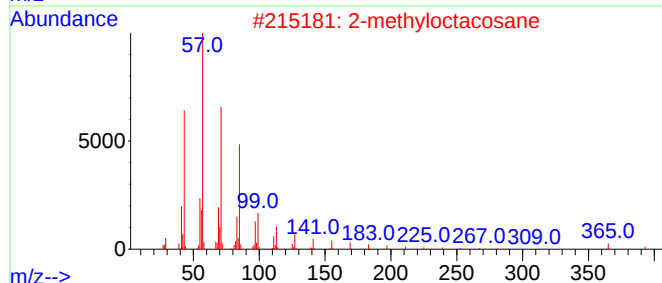
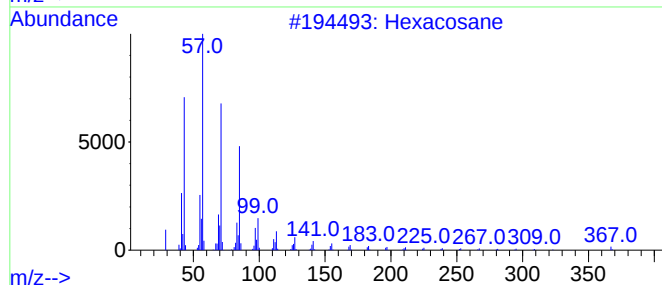
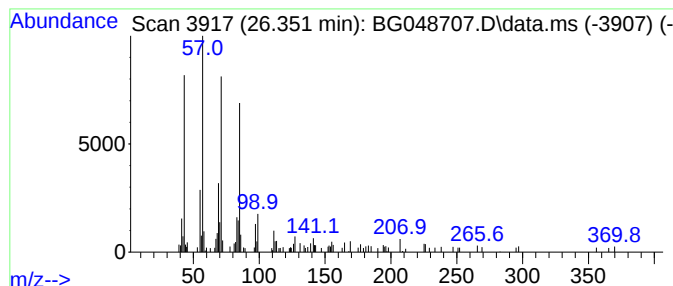
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.351	9.44 ng	191385	Perylene-d12	25.135

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			2-methyloctacosane	408	C29H60	1000376-72-8	87
3			Heptacosane	380	C27H56	000593-49-7	86
4			Tetracosane	338	C24H50	000646-31-1	86
5			Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

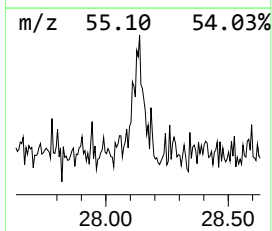
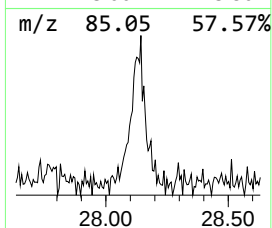
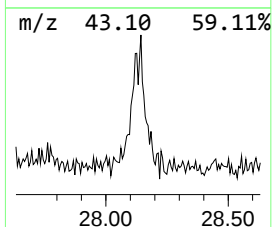
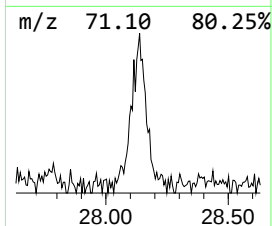
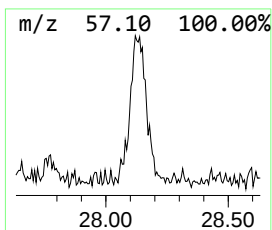
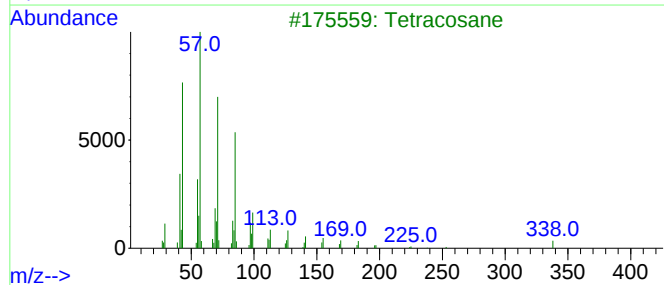
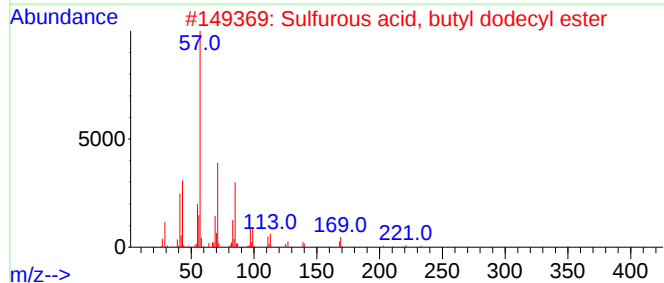
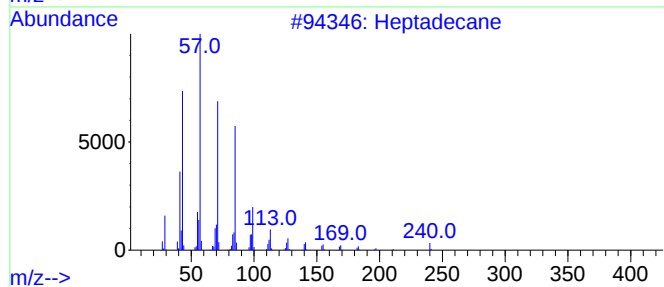
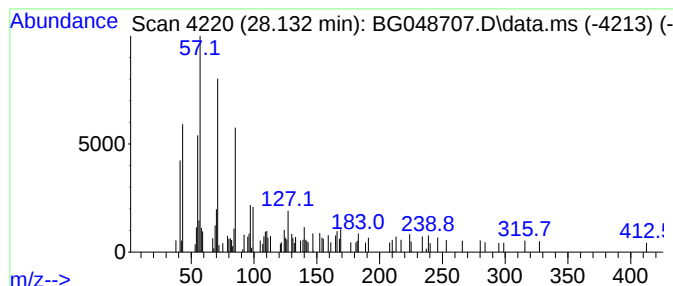
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Sulfurous acid, butyl dodec... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.132	7.17 ng	145348	Perylene-d12	25.135

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	86
2			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	64
3			Tetracosane	338	C24H50	000646-31-1	64
4			5,5-Diethylpentadecane	268	C19H40	1000360-41-0	59
5			2-methyltetracosane	352	C25H52	1000376-72-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041521\
Data File : BG048707.D
Acq On : 15 Apr 2021 13:40
Operator : CG/JU
Sample : M2015-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CL-02-041221

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,4,7,9-Tetrame...	13.619	6.9	ng	102038	3	14.736	295219	20.0
unknown15.846	15.846	4.5	ng	66353	3	14.736	295219	20.0
Pentadecanoic acid	18.372	5.8	ng	108722	4	17.491	373555	20.0
Benzamide, N-ac...	19.072	111.2	ng	2076650	4	17.491	373555	20.0
(3-Phenyloxiran...	19.213	135.4	ng	2529880	4	17.491	373555	20.0
2,2'-(Ethane-1,...	19.324	341.1	ng	6371510	4	17.491	373555	20.0
Morpholine, 4-p...	19.465	21.9	ng	409636	4	17.491	373555	20.0
Phthalic acid, ...	19.642	5.8	ng	108880	4	17.491	373555	20.0
Heptadecane	20.905	10.0	ng	186821	5	21.792	374198	20.0
Cycloeicosane	21.416	7.3	ng	136740	5	21.792	374198	20.0
Octacosane	21.704	9.4	ng	175573	5	21.792	374198	20.0
Hexacosane	22.621	13.5	ng	253155	5	21.792	374198	20.0
Heptadecane, 9-...	23.666	12.0	ng	242757	6	25.135	405388	20.0
Docosane, 5-butyl-	24.900	11.6	ng	235101	6	25.135	405388	20.0
2-methyloctacosane	26.351	9.4	ng	191385	6	25.135	405388	20.0
Sulfurous acid,...	28.132	7.2	ng	145348	6	25.135	405388	20.0