

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
 Data File : BF110521.D
 Acq On : 6 Nov 2018 18:02
 Operator : JU/SJ
 Sample : J5822-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-9

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.275	26	32	48	rVB	1642349	2267214	70.88%	9.003%
2	3.146	176	180	187	rBV	36189	60590	1.89%	0.241%
3	5.575	589	593	612	rBV	1148543	1120979	35.04%	4.451%
4	6.575	759	763	780	rBV2	749837	835812	26.13%	3.319%
5	6.722	784	788	791	rBV	2374992	2132610	66.67%	8.468%
6	6.951	824	827	833	rBV	747690	627422	19.61%	2.491%
7	7.110	850	854	857	rBV	2321339	2002256	62.59%	7.951%
8	7.522	918	924	944	rBV	1567117	1725652	53.95%	6.852%
9	8.233	1041	1045	1061	rBV	748786	864314	27.02%	3.432%
10	9.310	1223	1228	1231	rBV	3499980	3198852	100.00%	12.702%
11	9.704	1291	1295	1306	rVB	101605	115068	3.60%	0.457%
12	9.992	1341	1344	1357	rVB	951583	1022559	31.97%	4.060%
13	10.792	1475	1480	1493	rBV	1838549	1947683	60.89%	7.734%
14	11.486	1594	1598	1612	rBV	1082141	1065279	33.30%	4.230%
15	11.992	1680	1684	1698	rBV2	144606	187892	5.87%	0.746%
16	12.774	1814	1817	1826	rBV	182320	220554	6.89%	0.876%
17	13.074	1862	1868	1871	rBV	2934751	2999328	93.76%	11.910%
18	13.615	1956	1960	1968	rVB	28184	37169	1.16%	0.148%
19	13.945	2012	2016	2023	rBV	96801	118937	3.72%	0.472%
20	14.133	2044	2048	2057	rBV	615540	635151	19.86%	2.522%
21	14.257	2066	2069	2073	rVB	58226	53259	1.66%	0.211%
22	14.562	2118	2121	2128	rBV	108902	104905	3.28%	0.417%
23	14.880	2171	2175	2179	rBV	157655	154983	4.84%	0.615%
24	15.233	2230	2235	2240	rBV	171571	221285	6.92%	0.879%
25	15.627	2296	2302	2304	rBV	189277	262701	8.21%	1.043%
26	15.656	2304	2307	2319	rVB	338754	519446	16.24%	2.063%
27	16.080	2373	2379	2386	rBV	155208	239069	7.47%	0.949%
28	16.345	2419	2424	2429	rBV2	111913	175374	5.48%	0.696%
29	16.609	2462	2469	2475	rBV	83623	155754	4.87%	0.618%
30	17.233	2570	2575	2582	rVB3	28847	54234	1.70%	0.215%
31	17.592	2629	2636	2645	rVB6	25970	57089	1.78%	0.227%

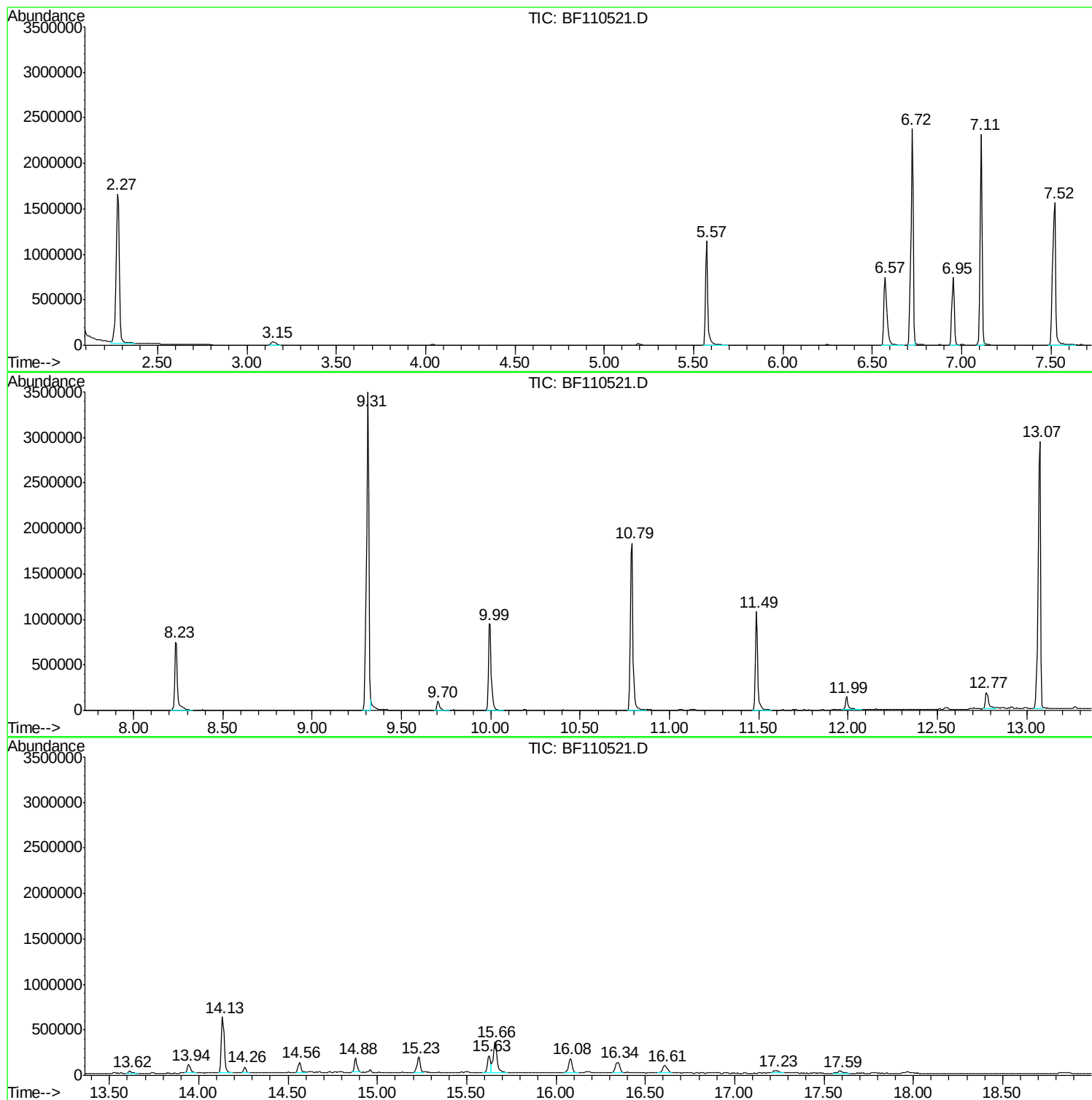
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BNA_F
ClientSampled :
MW-9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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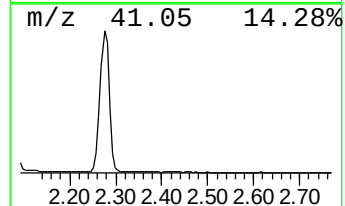
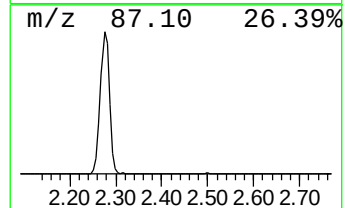
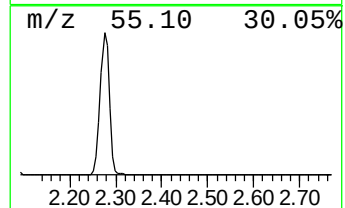
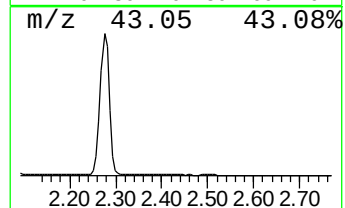
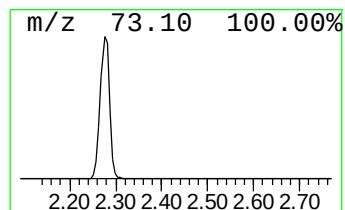
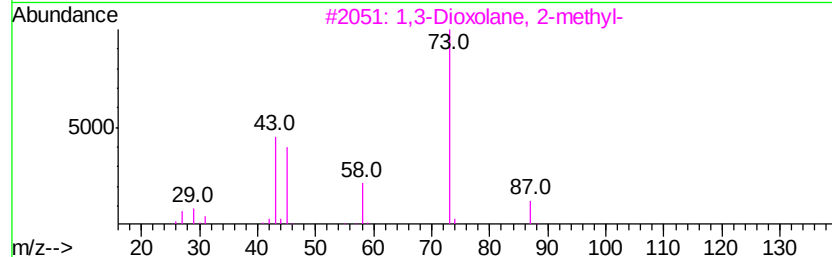
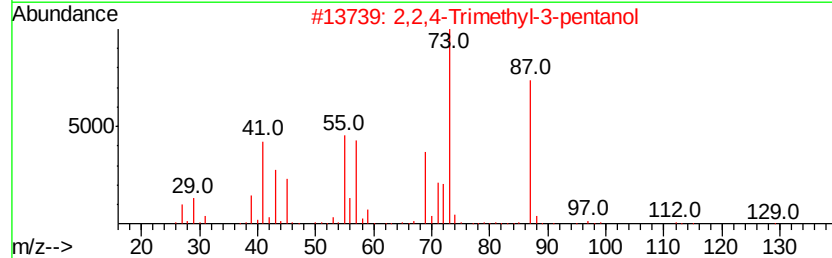
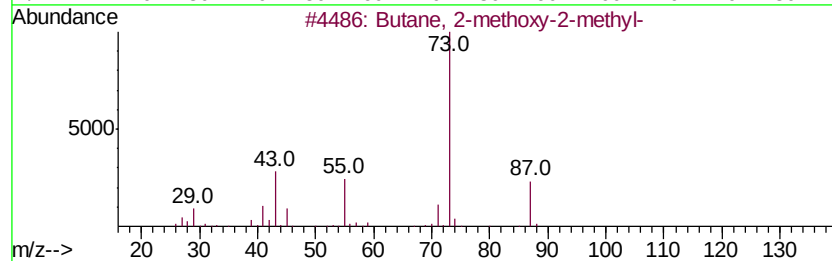
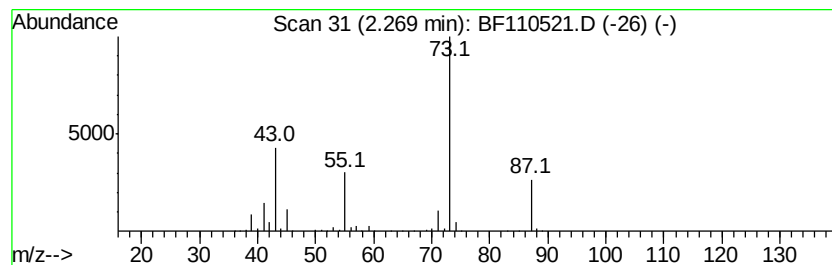
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	72.27 ng	2267210	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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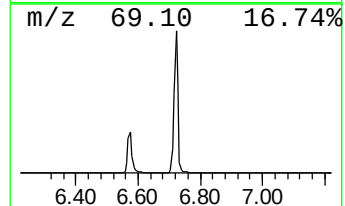
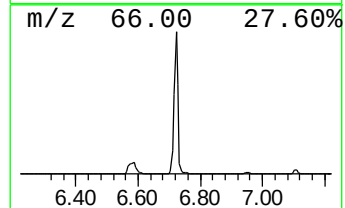
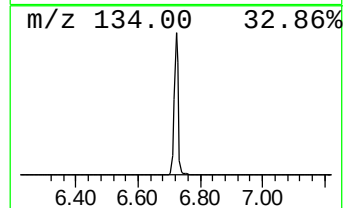
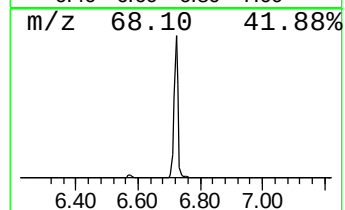
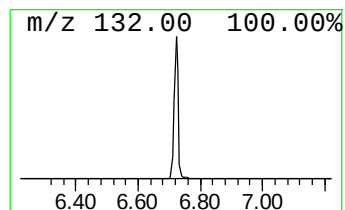
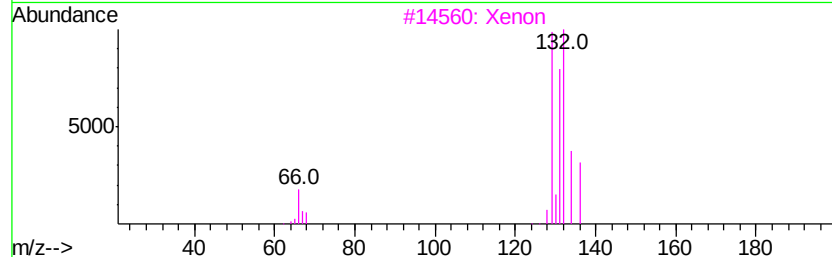
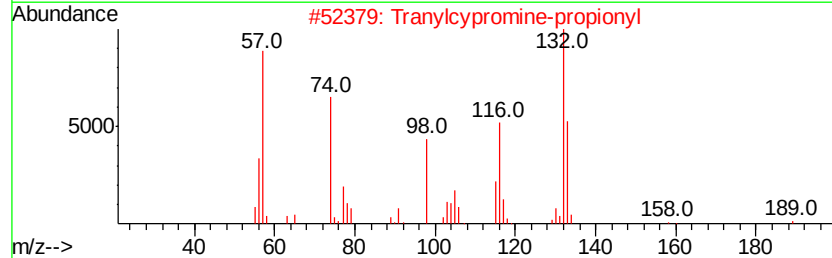
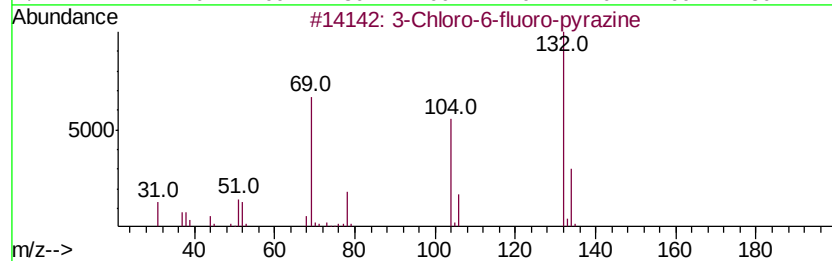
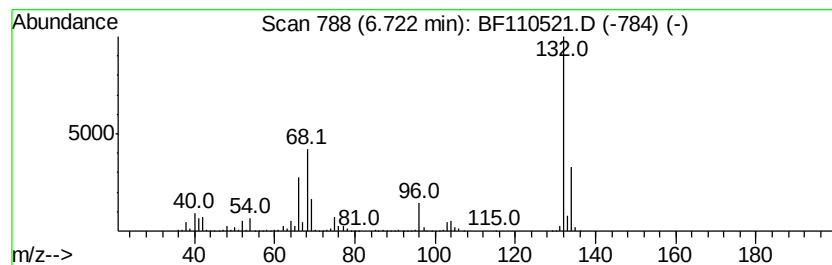
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Peak Number 2 unknown6.72 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	67.98 ng	2132610	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Xenon	132	Xe	007440-63-3	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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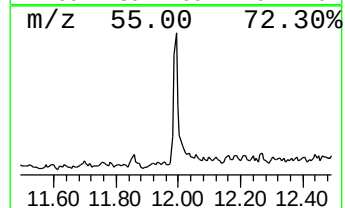
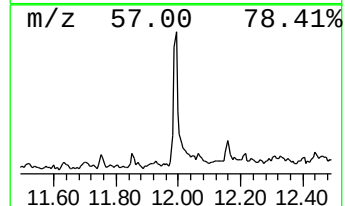
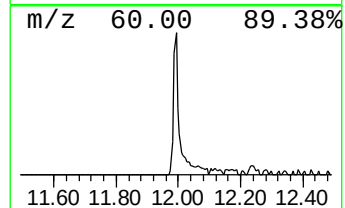
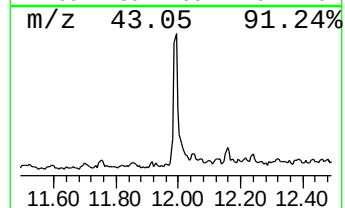
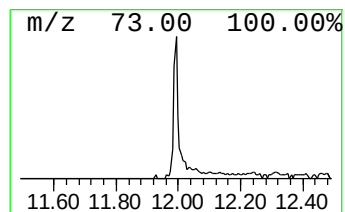
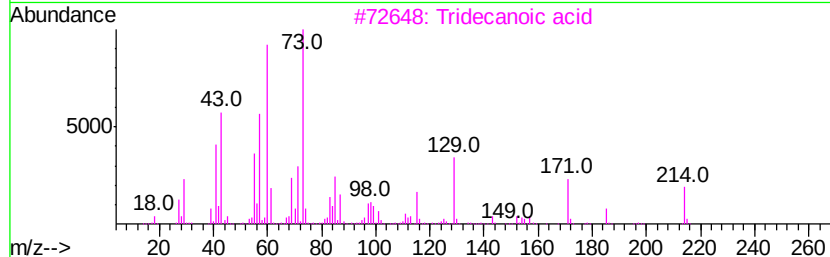
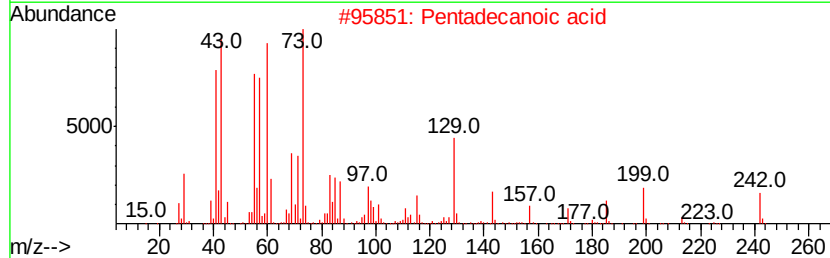
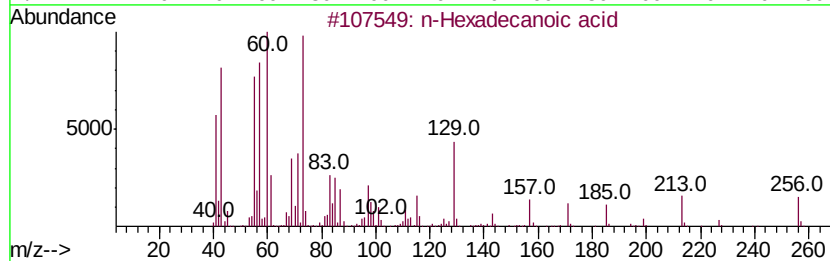
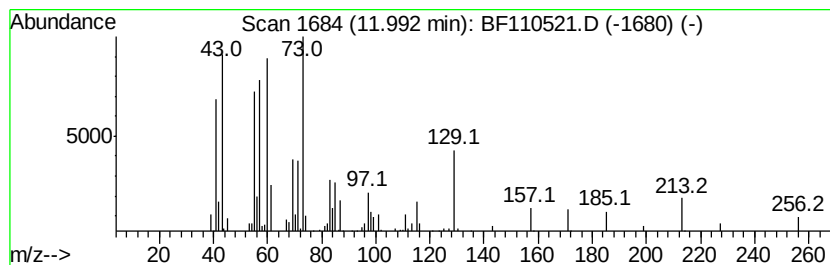
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	3.53 ng	187892	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tridecanoic acid	214	C13H26O2	000638-53-9	76
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



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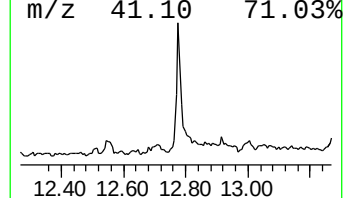
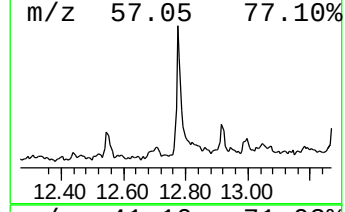
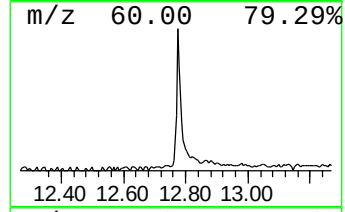
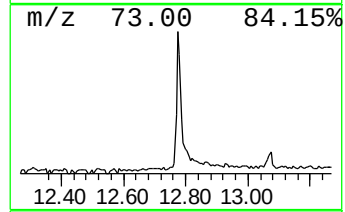
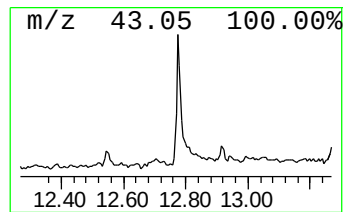
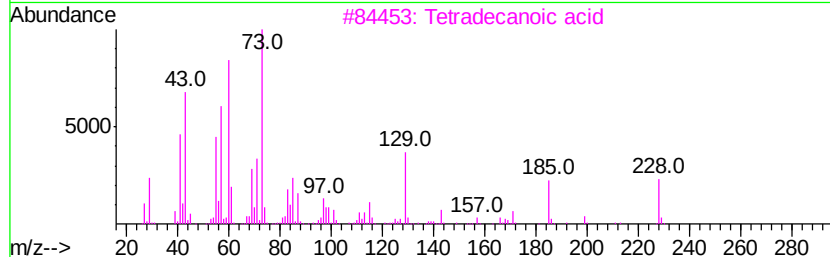
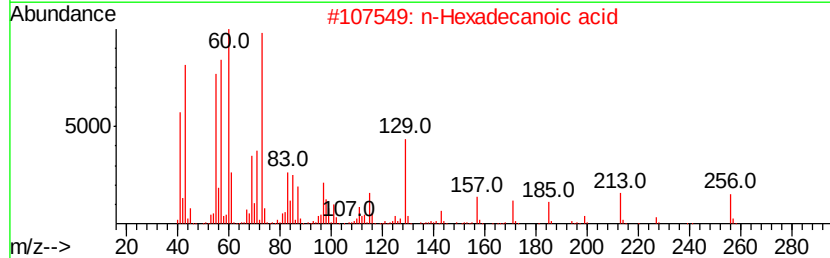
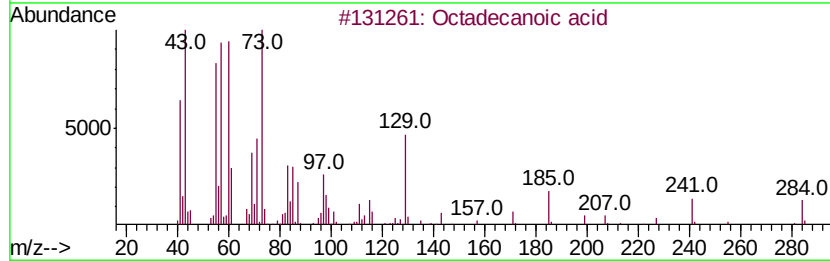
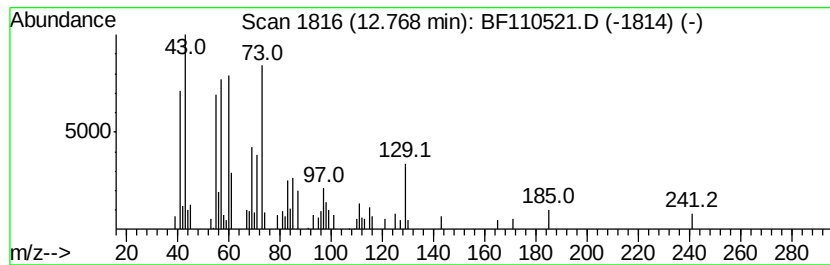
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	4.14 ng	220554	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5		Undecanoic acid	186	C11H22O2	000112-37-8	62



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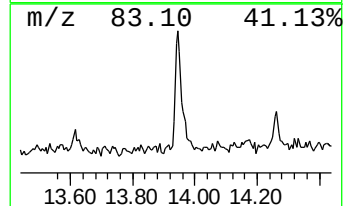
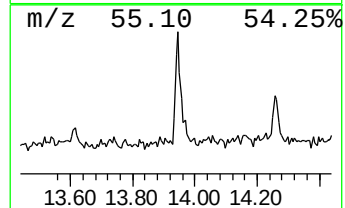
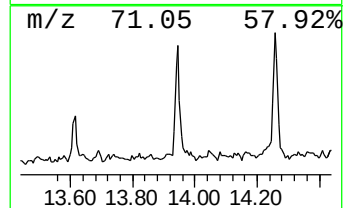
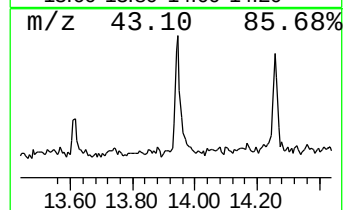
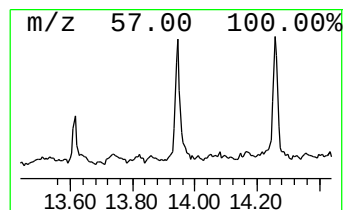
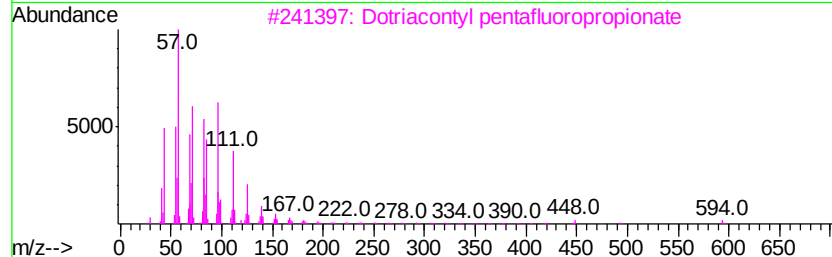
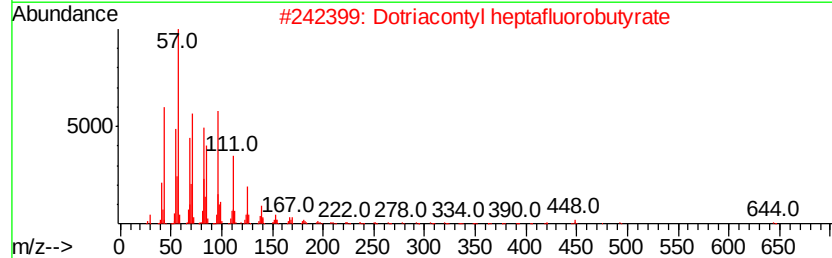
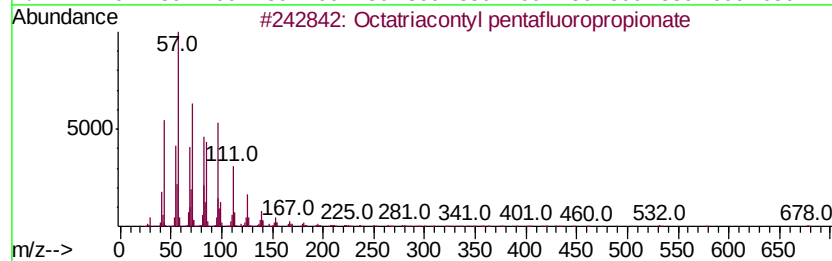
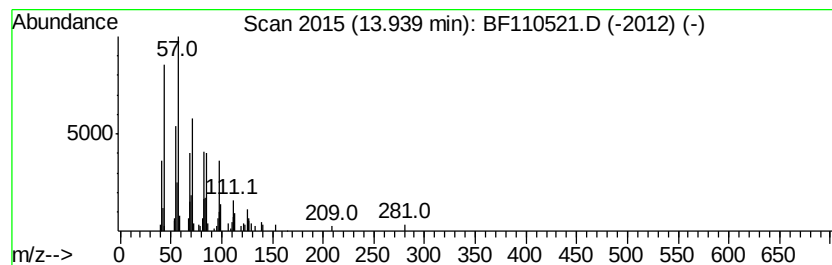
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Octatriacontyl pentafluorop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	3.75 ng	118937	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
2		Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	91
3		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	91
4		Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	91
5		Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	91



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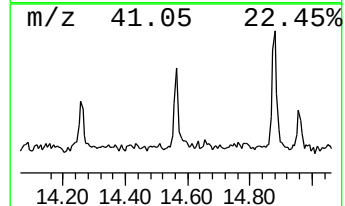
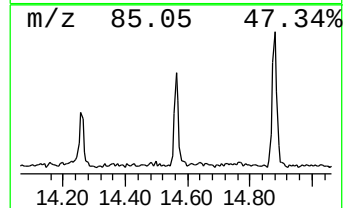
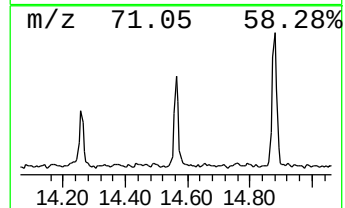
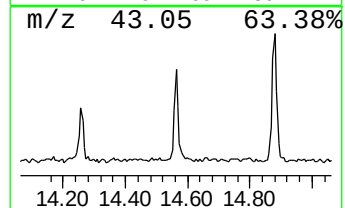
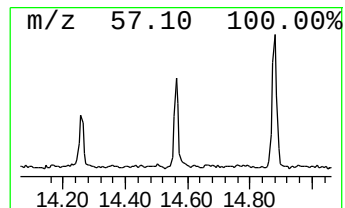
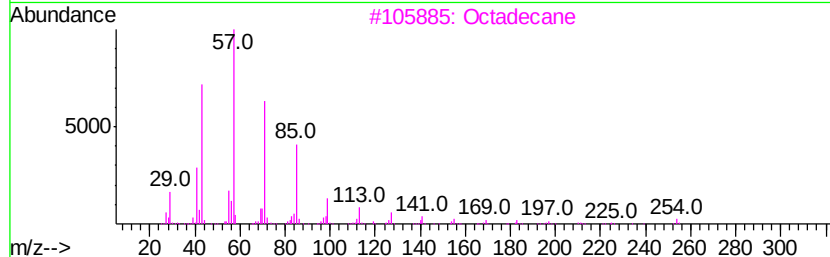
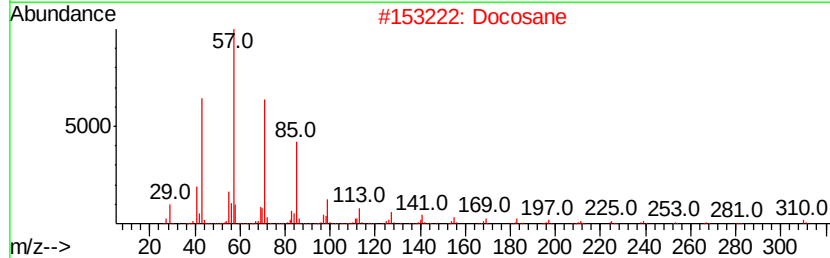
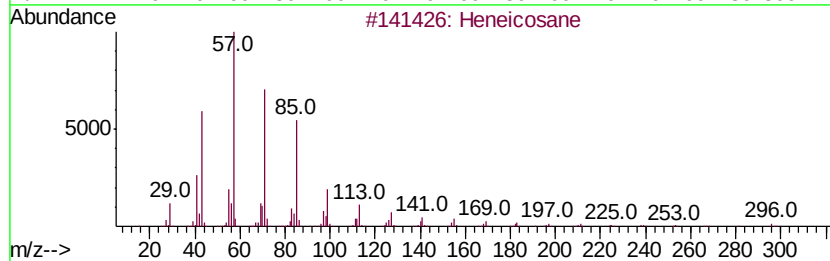
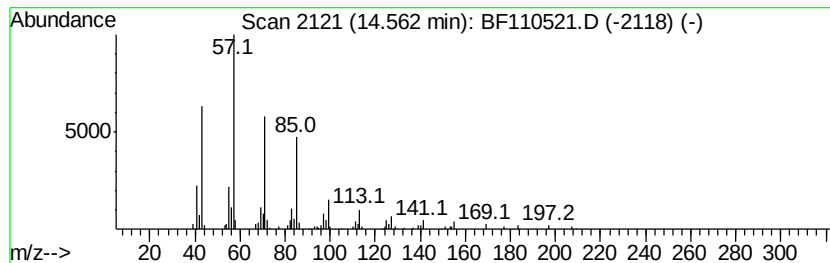
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	3.30 ng	104905	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Docosane	310	C22H46	000629-97-0	91
3		Octadecane	254	C18H38	000593-45-3	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
Operator : JU/SJ
Sample : J5822-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-9

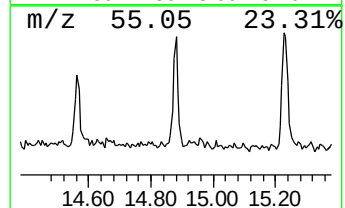
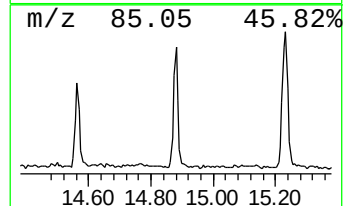
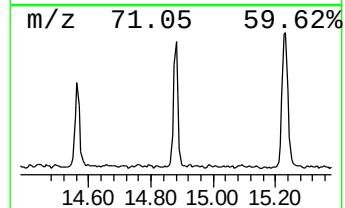
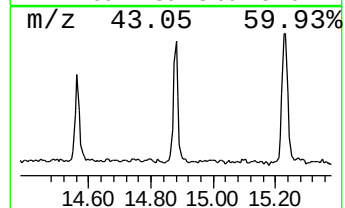
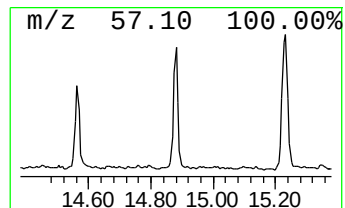
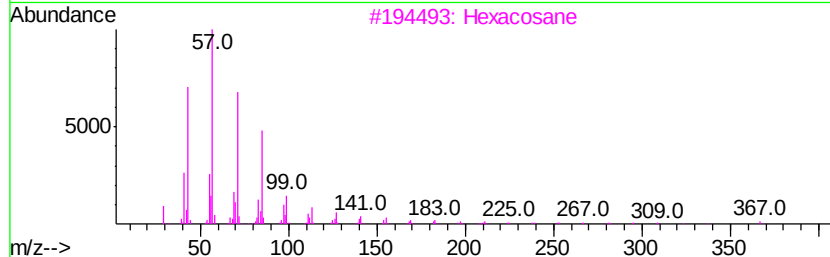
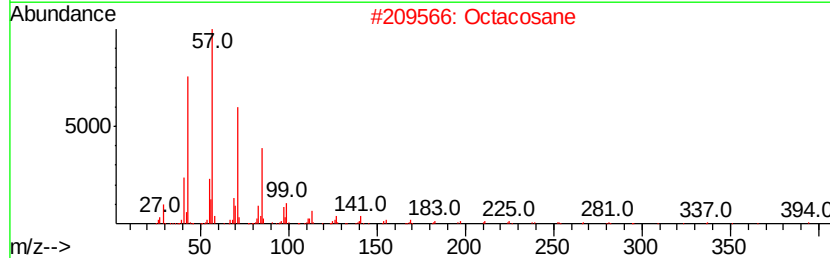
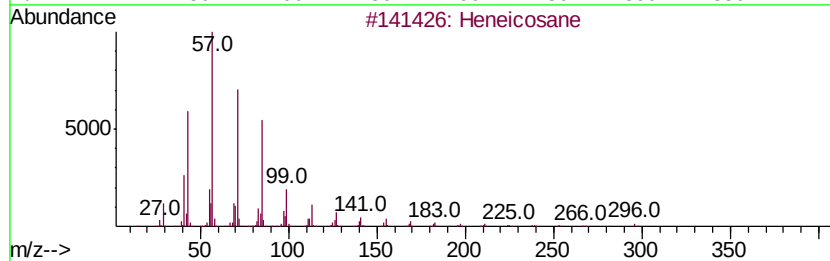
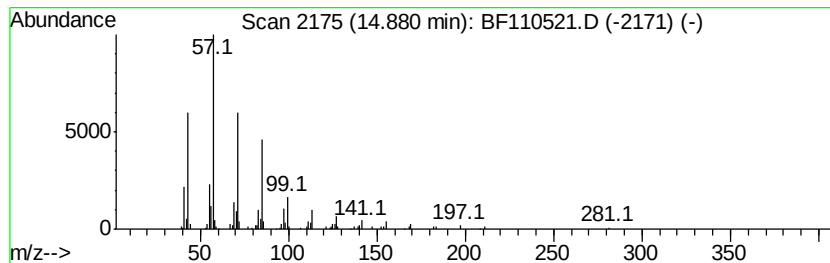
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	4.88 ng	154983	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Octacosane		394	C28H58	000630-02-4	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Pentacosane		352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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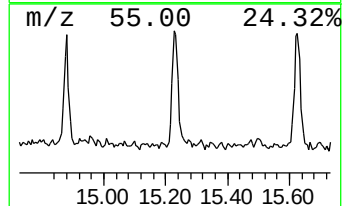
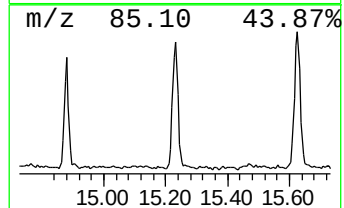
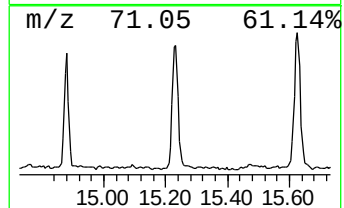
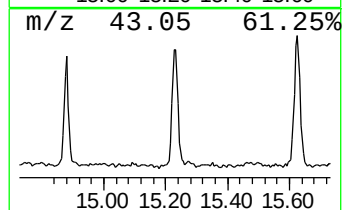
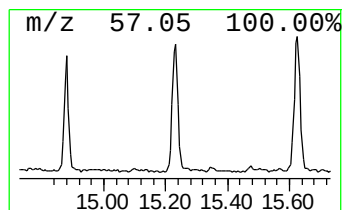
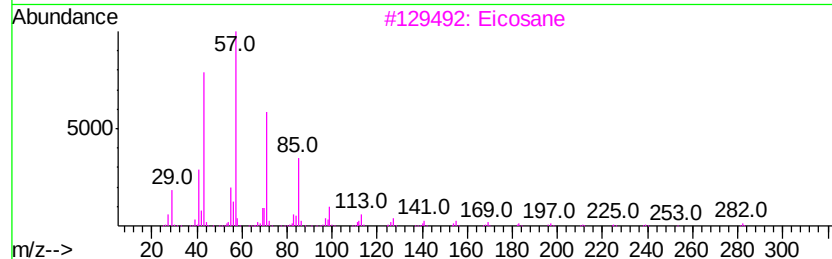
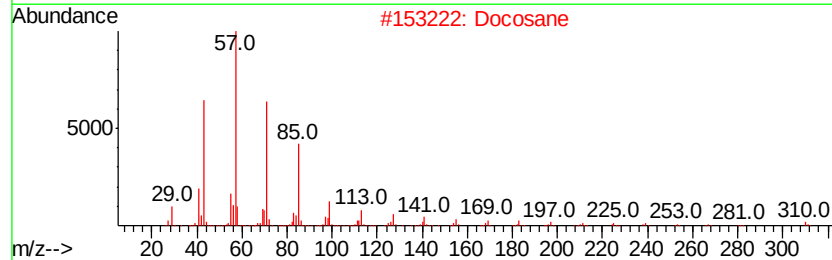
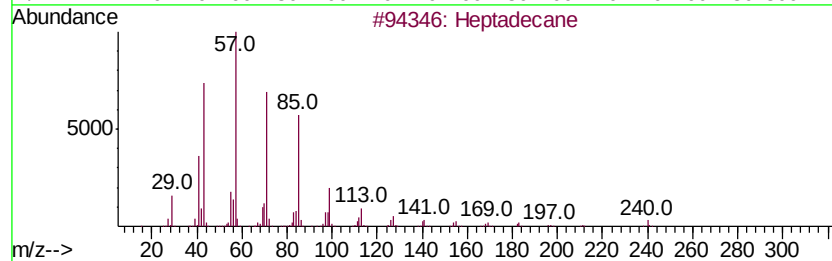
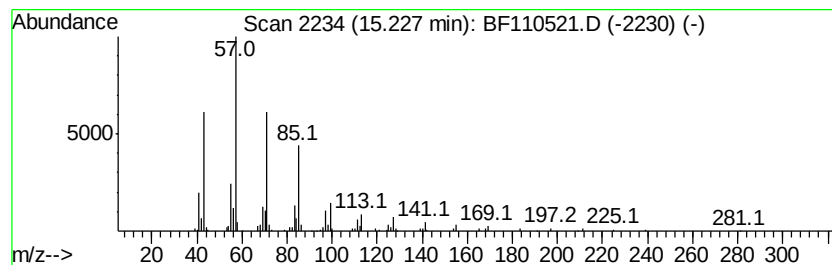
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.23	8.52 ng	221285	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Docosane	310	C22H46	000629-97-0	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
Operator : JU/SJ
Sample : J5822-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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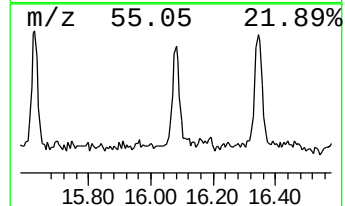
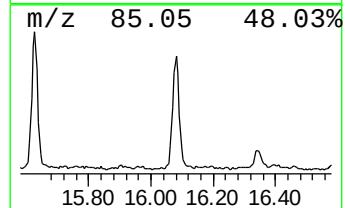
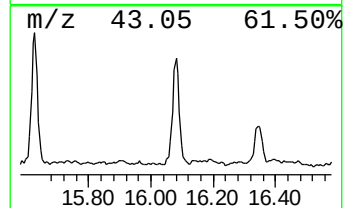
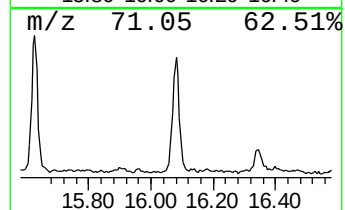
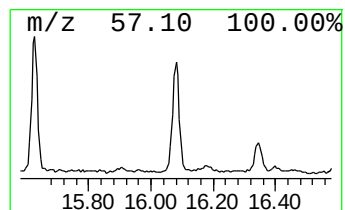
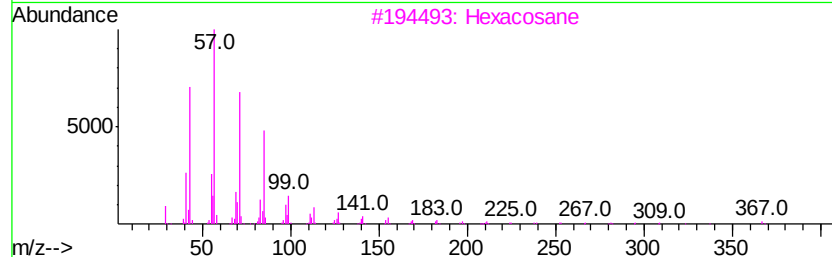
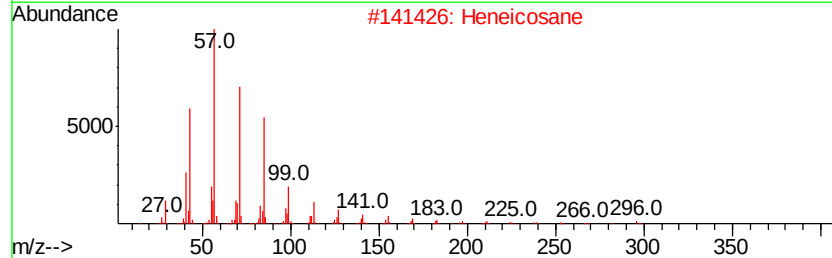
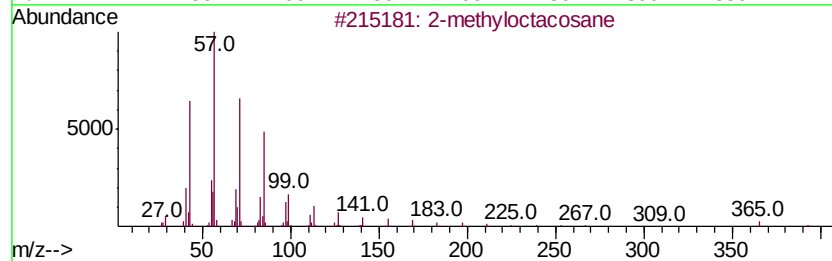
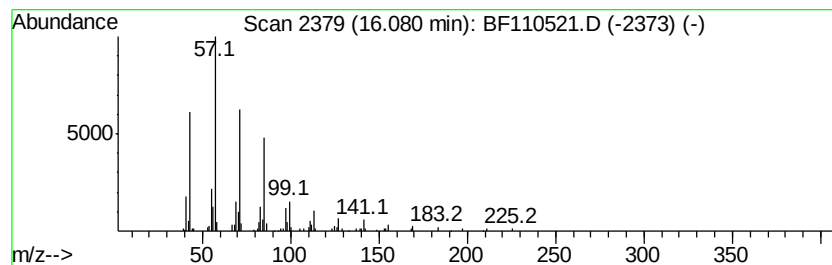
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-methyloctacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	9.20 ng	239069	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
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Operator : JU/SJ
Sample : J5822-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

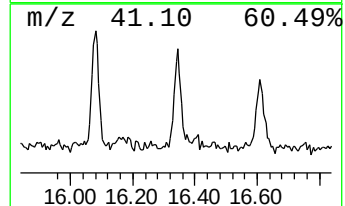
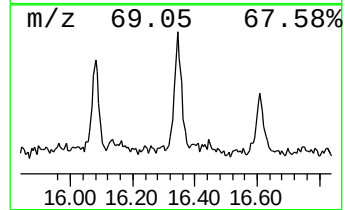
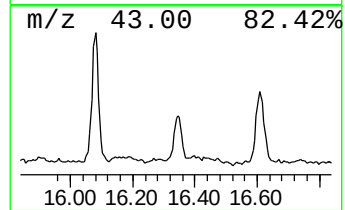
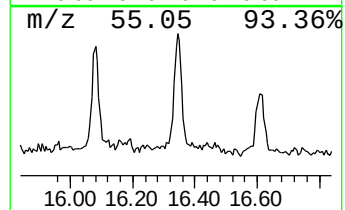
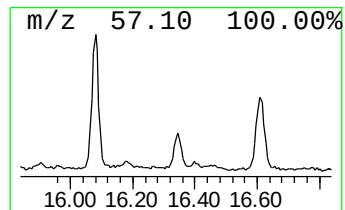
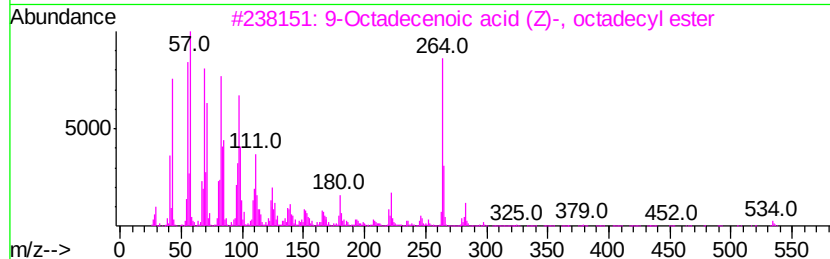
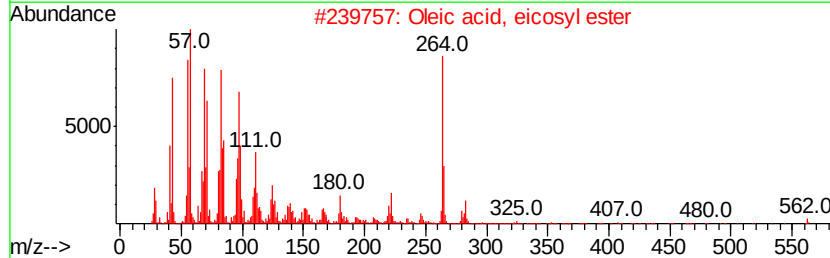
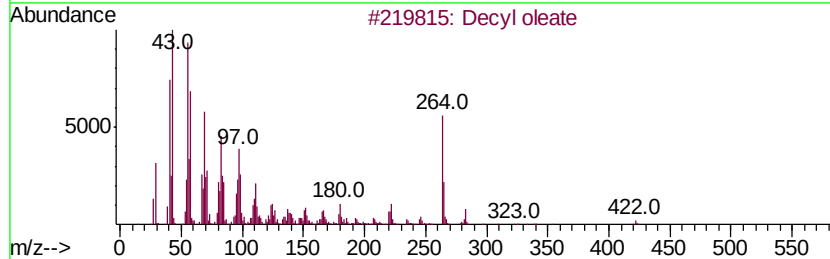
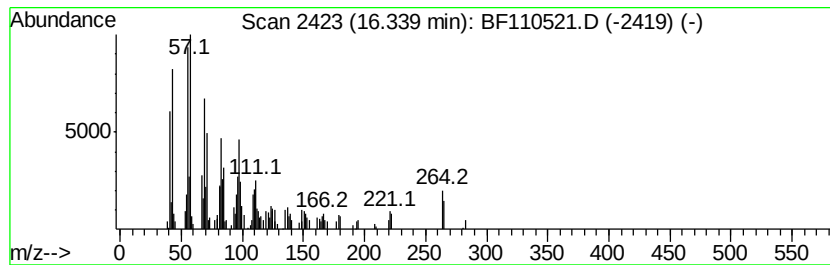
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Decyl oleate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	6.75 ng	175374	Perylene-d12	15.66
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Decyl oleate	422 C28H54O2	003687-46-5 91
2		Oleic acid, eicosyl ester	563 C38H74O2	022393-88-0 83
3		9-Octadecenoic acid (Z)-, octade...	535 C36H70O2	017673-49-3 83
4		10-Octadecenoic acid, methyl ester	296 C19H36O2	013481-95-3 76
5		cis-Vaccenic acid	282 C18H34O2	000506-17-2 74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
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Sample : J5822-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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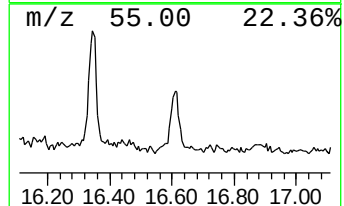
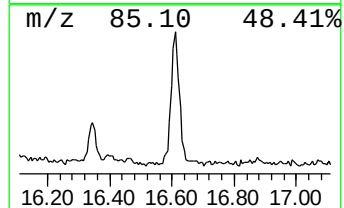
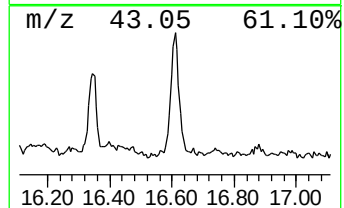
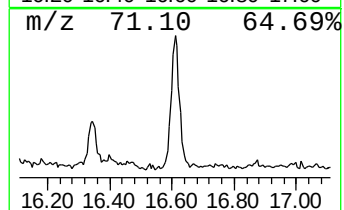
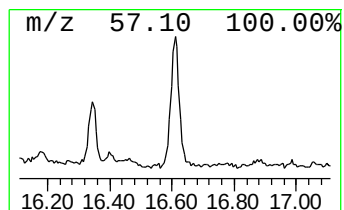
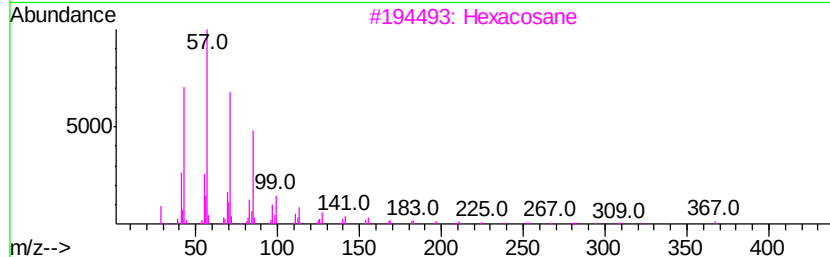
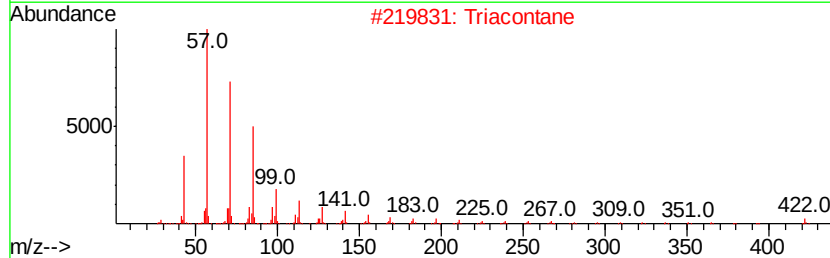
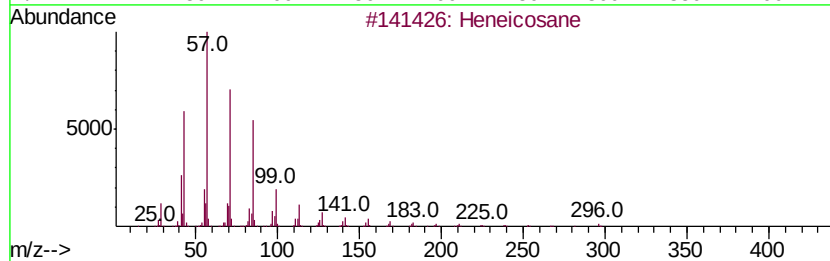
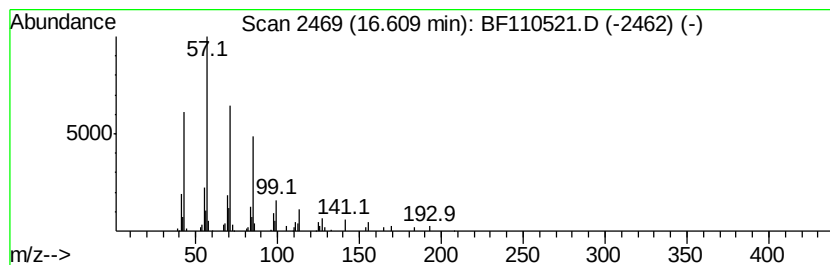
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Triacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.61	6.00 ng	155754	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Docosane	310	C22H46	000629-97-0	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
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Sample : J5822-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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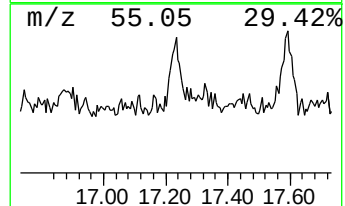
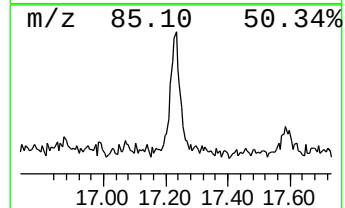
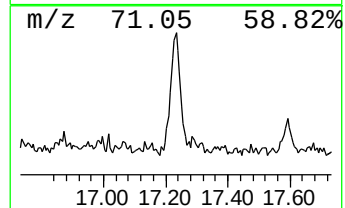
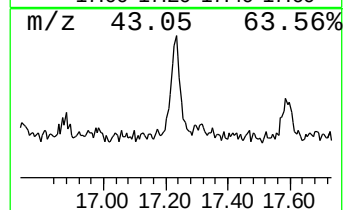
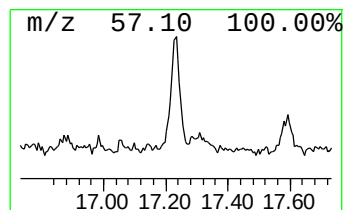
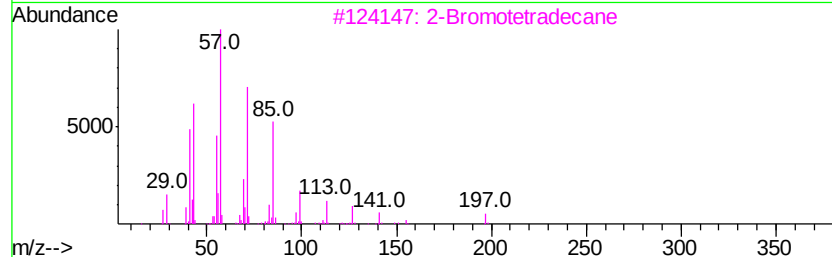
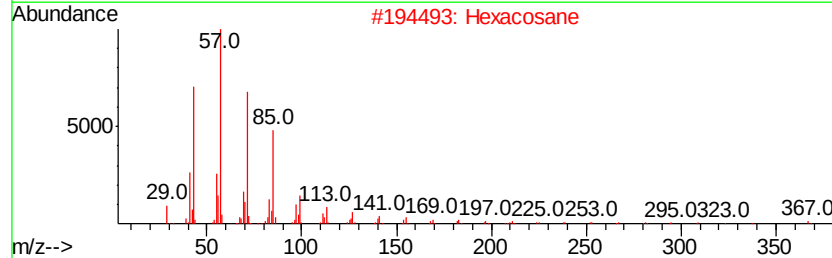
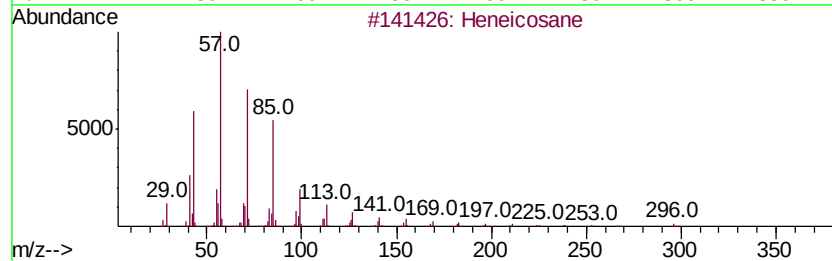
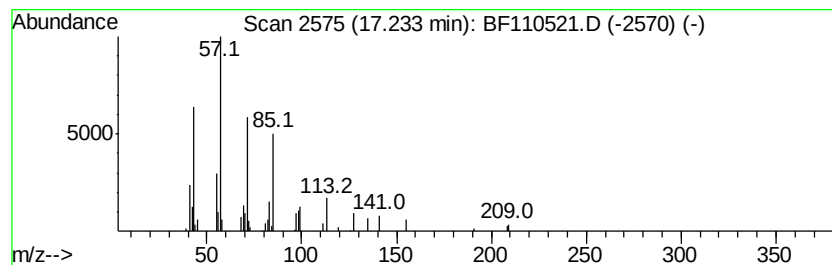
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-Bromotetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.23	2.09 ng	54234	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	80
2		Hexacosane	366	C26H54	000630-01-3	72
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	72
4		Pentacosane	352	C25H52	000629-99-2	72
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110521.D
Acq On : 6 Nov 2018 18:02
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ALS Vial : 7 Sample Multiplier: 1

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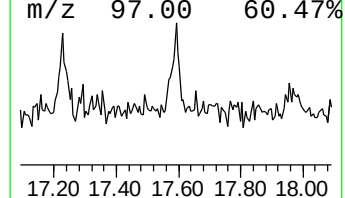
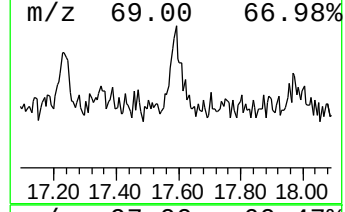
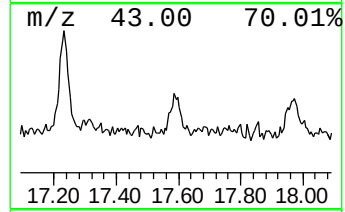
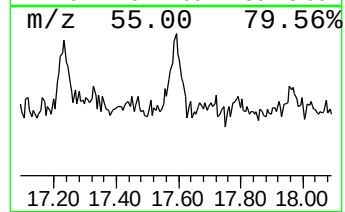
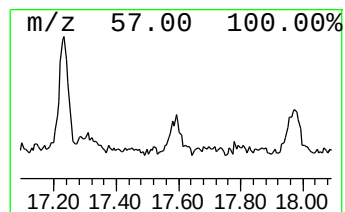
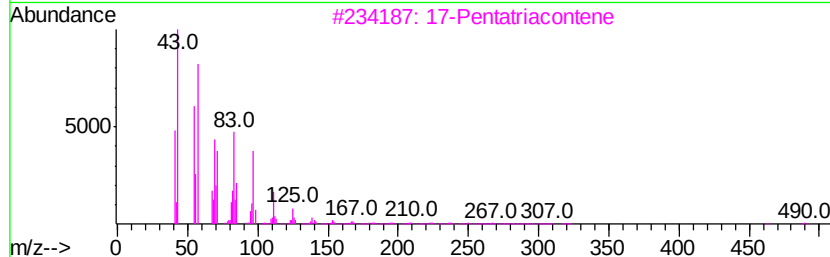
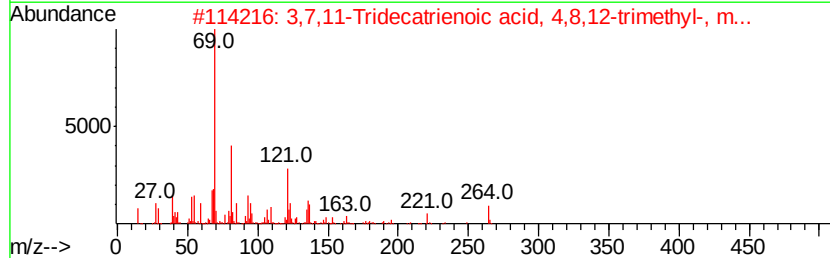
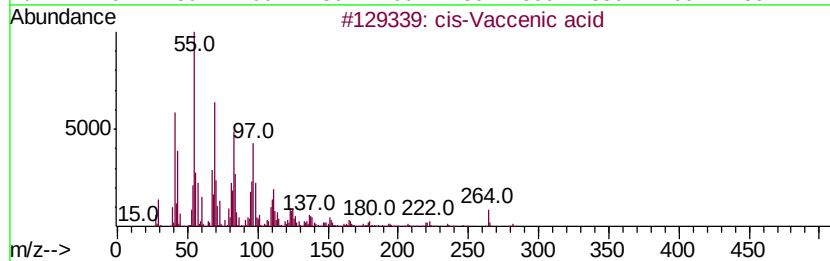
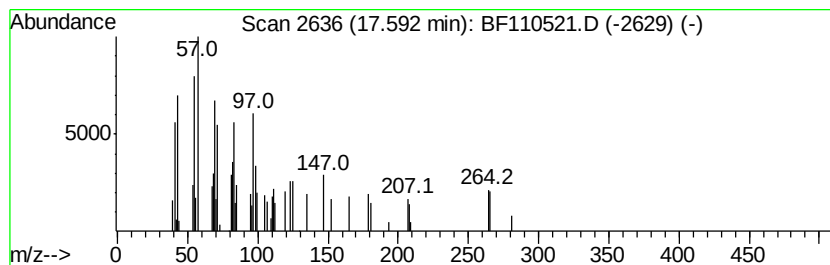
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.59 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.59	2.20 ng	57089	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	14
2		3,7,11-Tridecatrienoic acid, 4,8...	264	C17H28O2	036237-70-4	14
3		17-Pentatriacontene	491	C35H70	006971-40-0	11
4		2-Pentyne-1,4-diol, 1-(2-furanyl...	180	C10H12O3	1000351-33-1	11
5		6,10,14-Trimethyl-pentadecan-2-ol	270	C18H38O	1000143-49-8	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110618\
 Data File : BF110521.D
 Acq On : 6 Nov 2018 18:02
 Operator : JU/SJ
 Sample : J5822-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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
Butane, 2-methoxy...	2.27	72.3	ng	2267210	1	6.95	627422	20.0
unknown6.72	6.72	68.0	ng	2132610	1	6.95	627422	20.0
n-Hexadecanoic acid	11.99	3.5	ng	187892	4	11.49	1065280	20.0
Octadecanoic acid	12.77	4.1	ng	220554	4	11.49	1065280	20.0
Octatriacontyl pe...	13.94	3.8	ng	118937	5	14.13	635151	20.0
Heneicosane	14.56	3.3	ng	104905	5	14.13	635151	20.0
Octacosane	14.88	4.9	ng	154983	5	14.13	635151	20.0
Heptadecane	15.23	8.5	ng	221285	6	15.66	519446	20.0
2-methyloctacosane	16.08	9.2	ng	239069	6	15.66	519446	20.0
Decyl oleate	16.34	6.8	ng	175374	6	15.66	519446	20.0
Triacontane	16.61	6.0	ng	155754	6	15.66	519446	20.0
2-Bromotetradecane	17.23	2.1	ng	54234	6	15.66	519446	20.0
unknown17.59	17.59	2.2	ng	57089	6	15.66	519446	20.0