

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
 Data File : BF107487.D
 Acq On : 18 Jul 2018 19:33
 Operator : JU/SJ
 Sample : J4035-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-2

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-BF071618.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	5.231	487	492	500	rVB	336172	466360	10.96%	1.214%
2	5.619	552	558	561	rBV	3401047	3631931	85.35%	9.457%
3	6.613	721	727	730	rBV	3326770	3852582	90.53%	10.032%
4	6.760	746	752	755	rBV	3729898	3923231	92.19%	10.216%
5	6.801	756	759	764	rBV2	52381	53291	1.25%	0.139%
6	6.984	785	790	793	rBV	1374395	1101002	25.87%	2.867%
7	7.142	812	817	820	rVB	3371967	2983482	70.11%	7.769%
8	7.548	878	886	889	rBV	2564970	2535695	59.59%	6.603%
9	7.613	892	897	905	rVV6	38652	46653	1.10%	0.121%
10	8.266	1003	1008	1012	rBV	1514324	1293101	30.39%	3.367%
11	9.342	1185	1191	1194	rBV	4348690	4255528	100.00%	11.081%
12	9.442	1205	1208	1212	rVB4	61960	60331	1.42%	0.157%
13	9.730	1253	1257	1261	rBV	692464	554893	13.04%	1.445%
14	9.972	1294	1298	1302	rVB	263794	241251	5.67%	0.628%
15	10.030	1304	1308	1312	rVB	1542509	1431445	33.64%	3.727%
16	10.442	1375	1378	1381	rVB	73547	65300	1.53%	0.170%
17	10.819	1438	1442	1446	rBV	2382006	2267211	53.28%	5.904%
18	10.913	1453	1458	1463	rVB5	61176	95694	2.25%	0.249%
19	11.195	1504	1506	1512	rVB	79716	73015	1.72%	0.190%
20	11.366	1531	1535	1538	rBV4	46843	44525	1.05%	0.116%
21	11.519	1557	1561	1565	rBV	1322487	1289197	30.29%	3.357%
22	11.724	1592	1596	1598	rBV5	53952	51217	1.20%	0.133%
23	12.548	1732	1736	1739	rBV2	346778	330933	7.78%	0.862%
24	12.589	1741	1743	1748	rVB4	79638	91516	2.15%	0.238%
25	12.907	1795	1797	1801	rVB5	61898	54082	1.27%	0.141%
26	12.965	1804	1807	1808	rVV	132170	115889	2.72%	0.302%
27	12.983	1808	1810	1818	rVB3	148131	151386	3.56%	0.394%
28	13.107	1824	1831	1835	rBV	3273838	3208846	75.40%	8.356%
29	13.295	1859	1863	1865	rBV2	155993	143519	3.37%	0.374%
30	13.318	1865	1867	1870	rVB	198780	163148	3.83%	0.425%
31	13.618	1915	1918	1922	rVB5	48076	57261	1.35%	0.149%
32	13.665	1922	1926	1928	rBV	216072	221710	5.21%	0.577%
33	13.989	1977	1981	1985	rBV	396434	440628	10.35%	1.147%
34	14.171	2008	2012	2015	rBV	1099077	998794	23.47%	2.601%

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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

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35	14.307	2032	2035	2039	rBV	238689	256718	6.03%	0.668%
36	14.536	2072	2074	2079	rBV6	57215	76681	1.80%	0.200%
37	14.618	2085	2088	2093	rBV	217359	212241	4.99%	0.553%
38	14.942	2140	2143	2147	rVB	195595	174128	4.09%	0.453%
39	15.018	2154	2156	2160	rVB2	121468	135465	3.18%	0.353%
40	15.301	2201	2204	2207	rVB4	178629	161581	3.80%	0.421%
41	15.712	2269	2274	2279	rVB	706460	929871	21.85%	2.421%
42	16.183	2349	2354	2358	rVB8	98493	162604	3.82%	0.423%

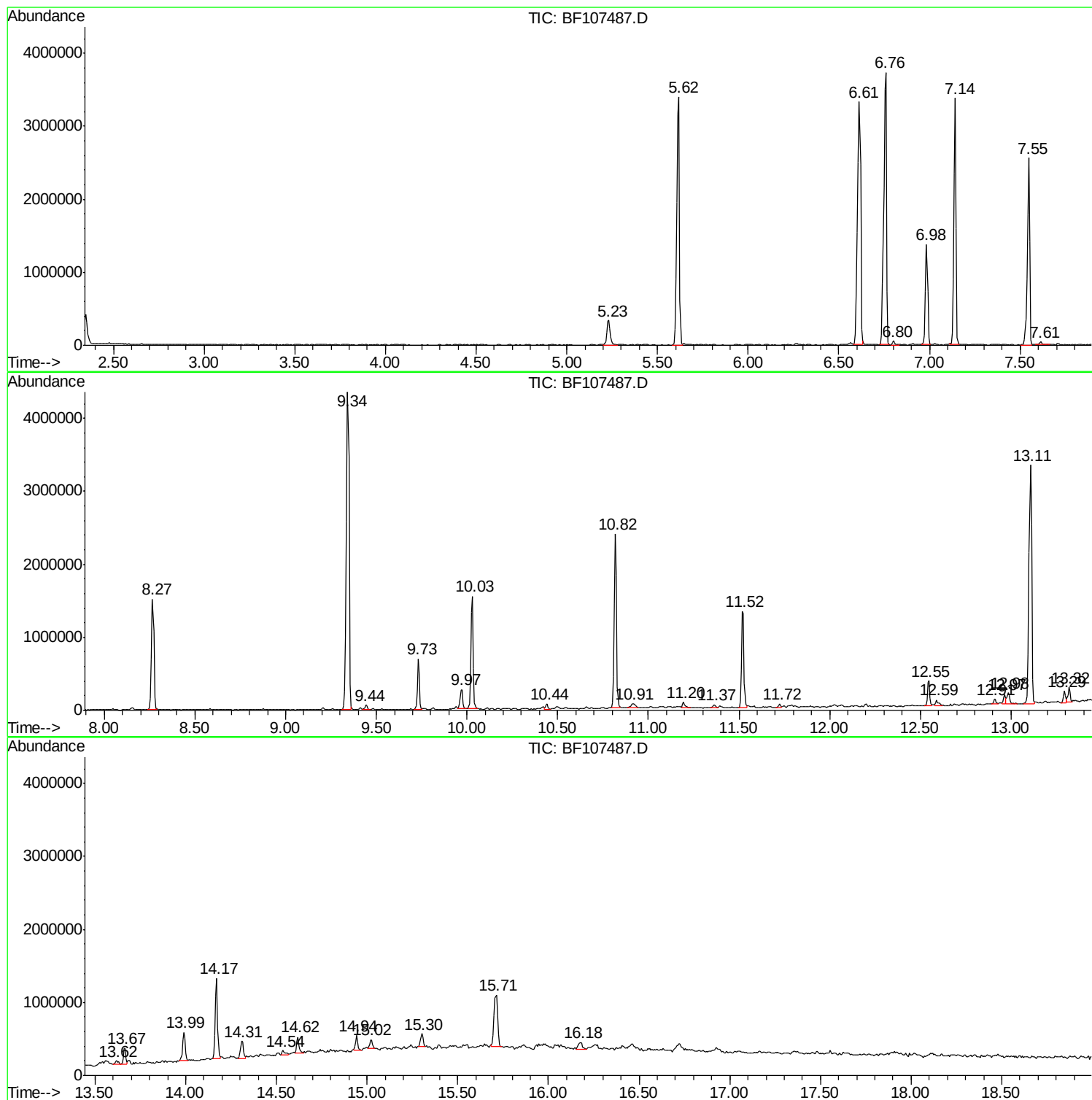
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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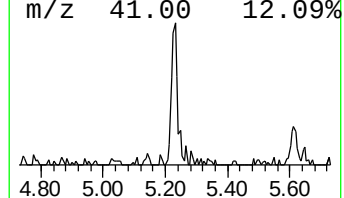
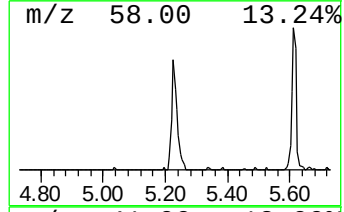
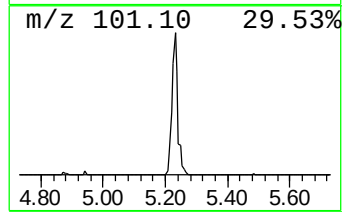
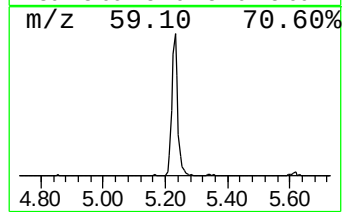
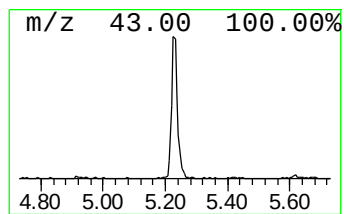
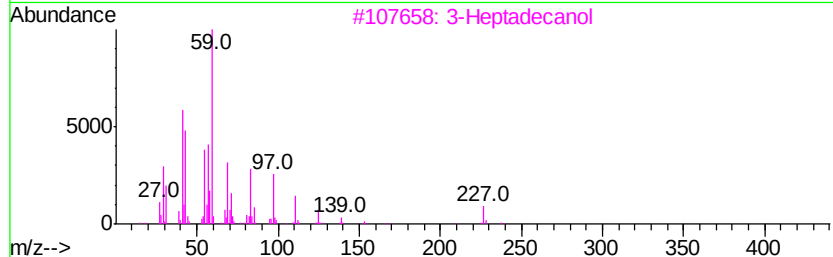
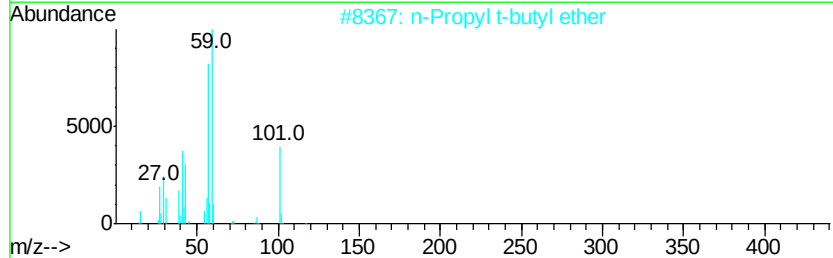
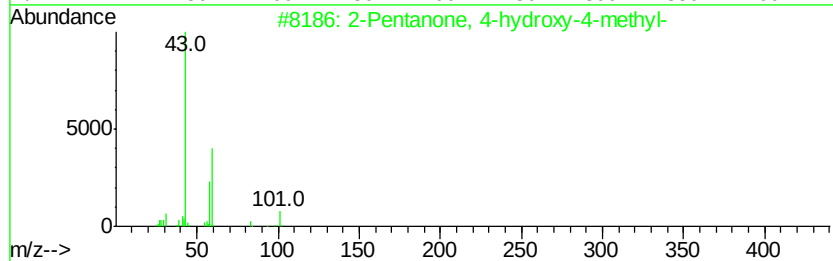
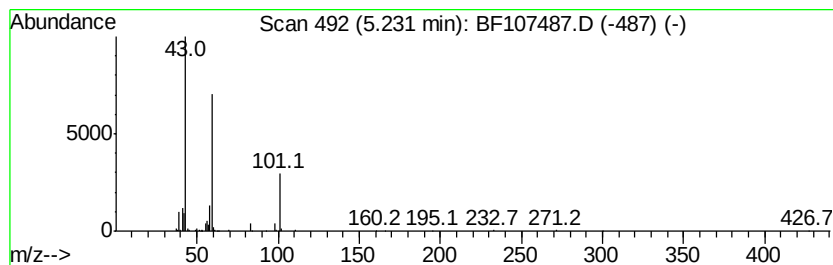
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	8.47 ng	466360	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	36
3			3-Heptadecanol	256	C17H36O	084534-30-5	23
4			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12



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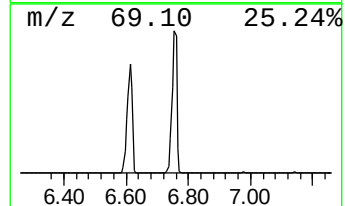
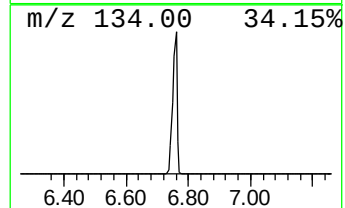
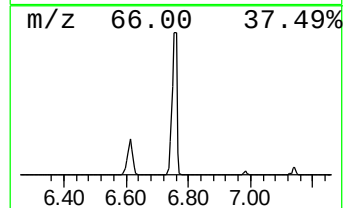
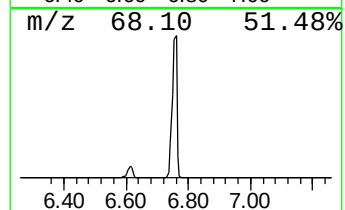
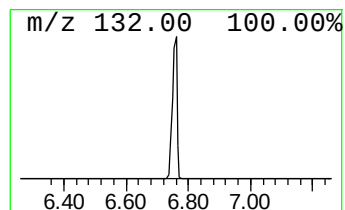
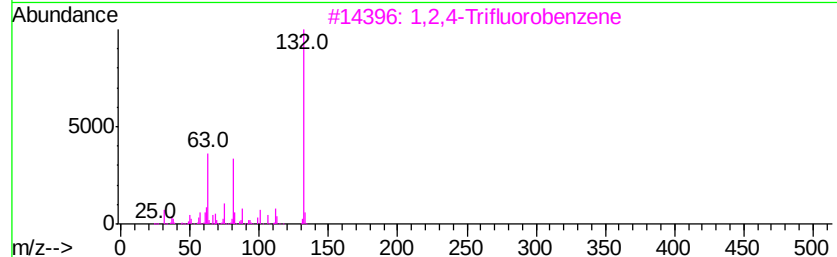
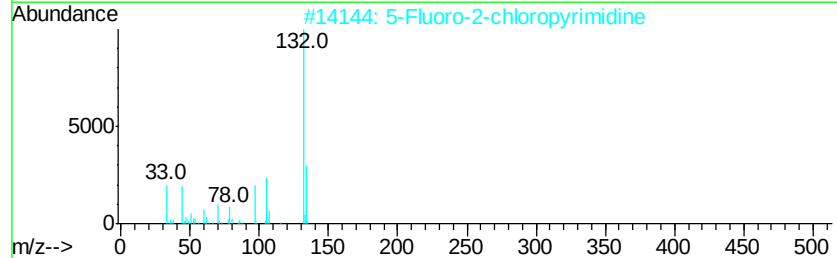
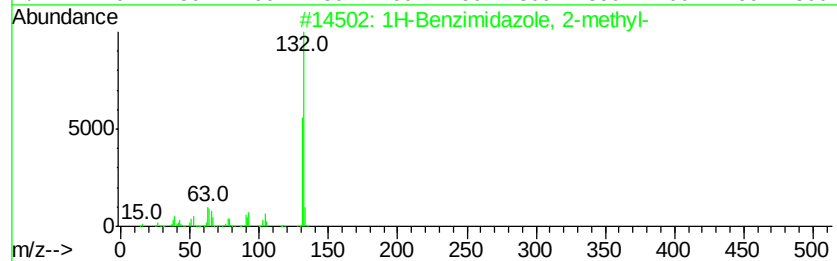
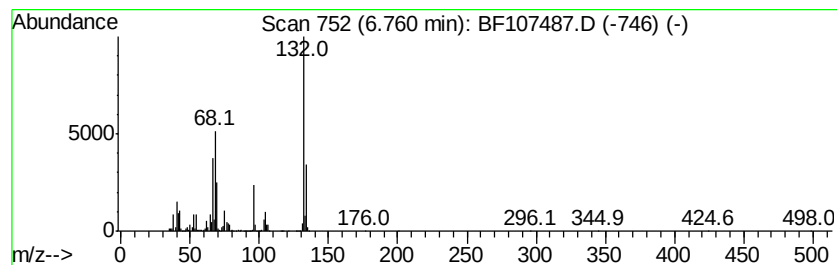
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Peak Number 2 unknown6.76 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	71.27 ng	3923230	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	14
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	14
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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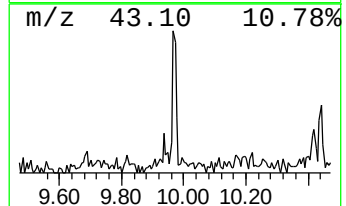
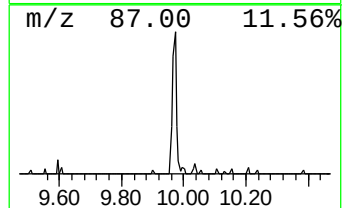
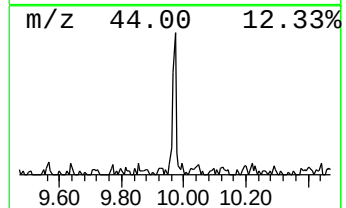
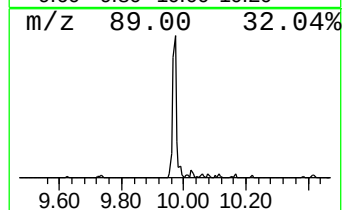
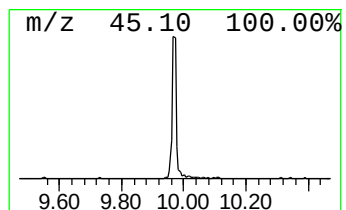
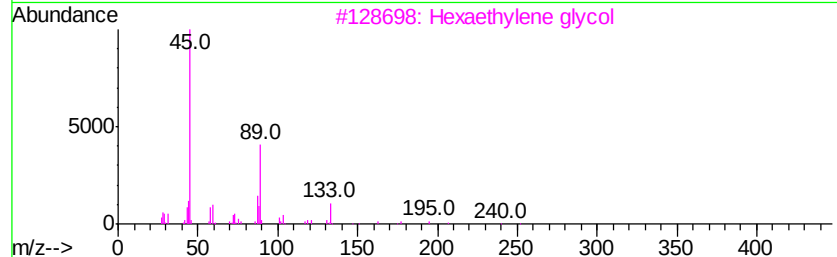
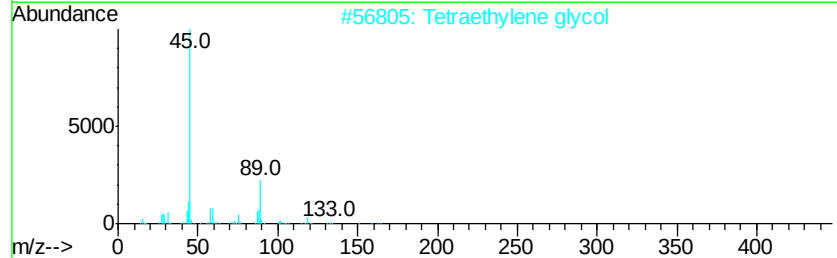
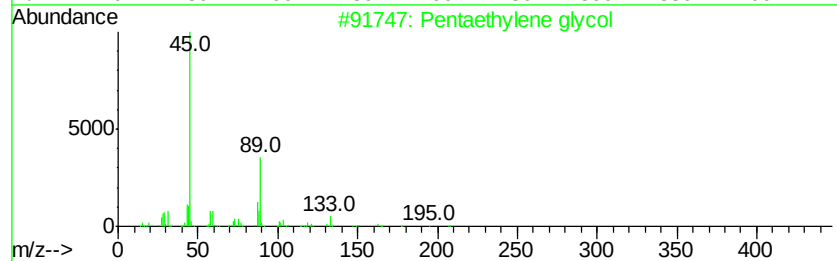
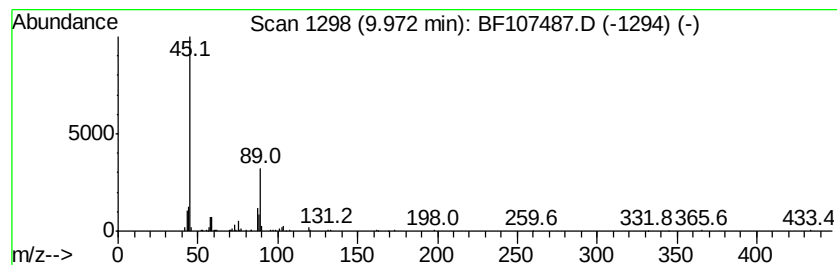
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentaethylene glycol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	3.37 ng	241251	Acenaphthene-d10	10.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentaethylene glycol	238	C10H22O6	004792-15-8	90	
2	Tetraethylene glycol	194	C8H18O5	000112-60-7	90	
3	Hexaethylene glycol	282	C12H26O7	002615-15-8	90	
4	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	78	
5	Ethanol, 2-[2-(2-methoxyethoxy)ethoxy]ethoxy	164	C7H16O4	000112-35-6	78	



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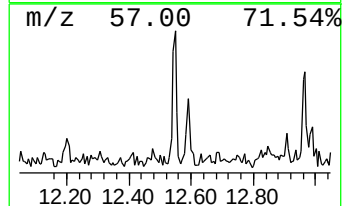
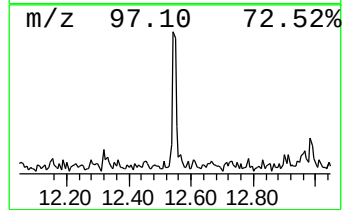
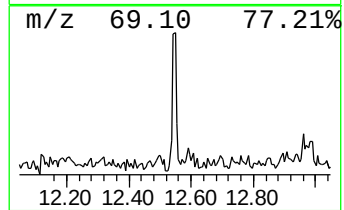
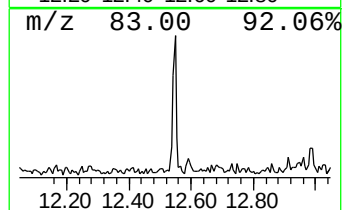
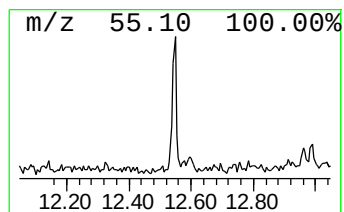
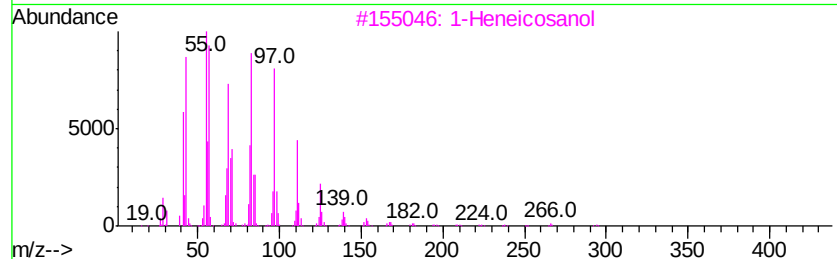
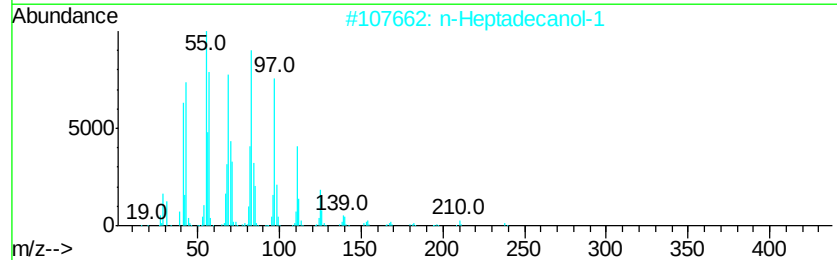
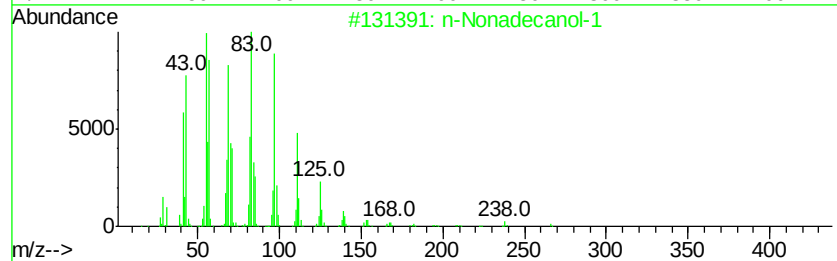
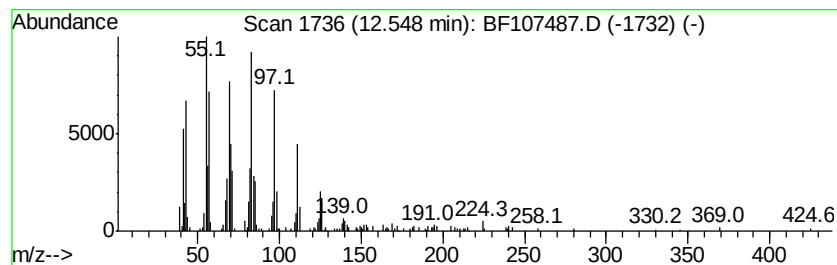
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 n-Nonadecanol-1 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	5.13 ng	330933	Phenanthrene-d10	11.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	n-Heptadecanol-1	256	C17H36O	001454-85-9	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	3-Heptadecene, (Z)-	238	C17H34	1000141-67-3	93
5	Cyclohexadecane	224	C16H32	000295-65-8	91



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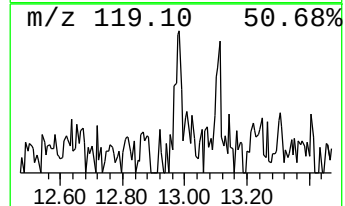
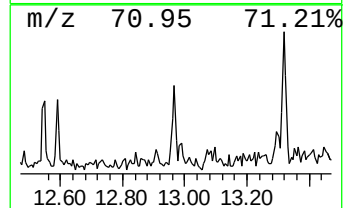
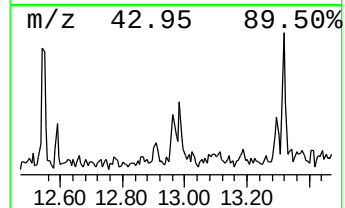
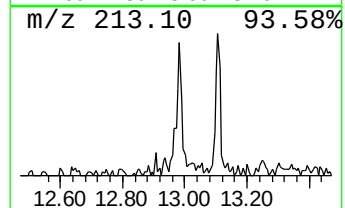
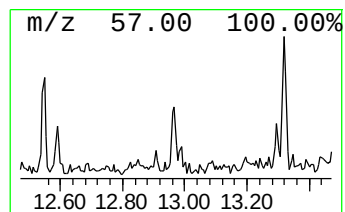
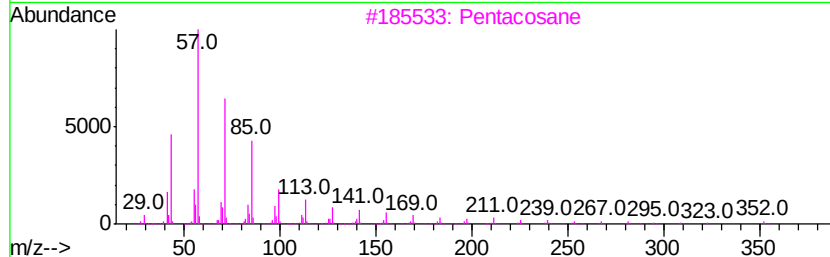
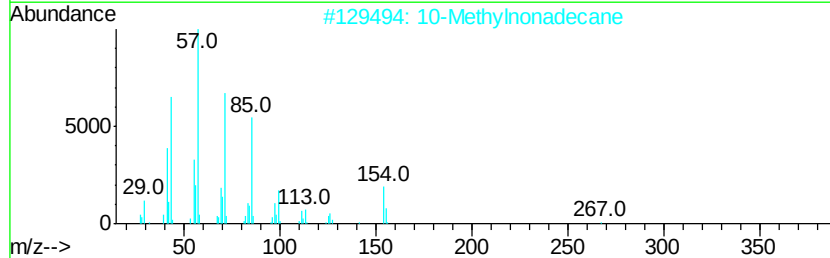
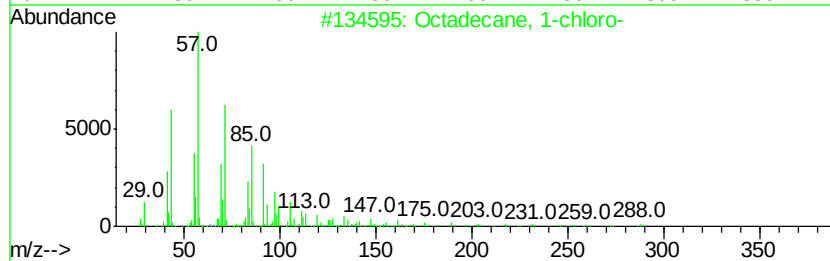
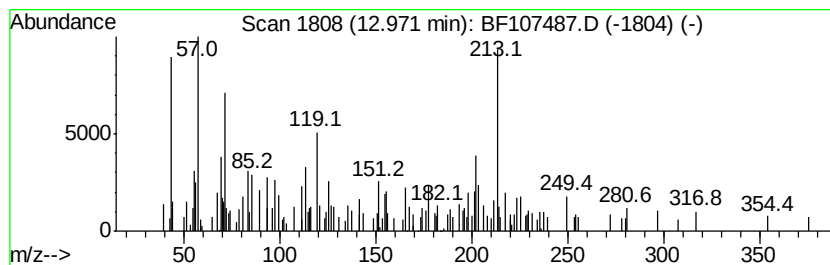
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Peak Number 6 Octadecane, 1-chloro- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.97	2.32 ng	115889	Chrysene-d12	14.17		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	70
2		10-Methylnonadecane	282	C20H42	056862-62-5	64
3		Pentacosane	352	C25H52	000629-99-2	62
4		Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	58
5		Heptadecane	240	C17H36	000629-78-7	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
Operator : JU/SJ
Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-2

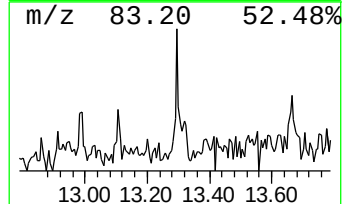
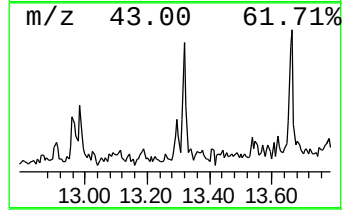
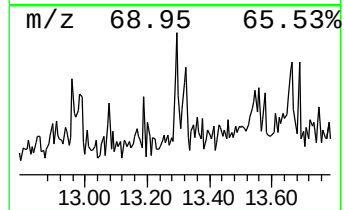
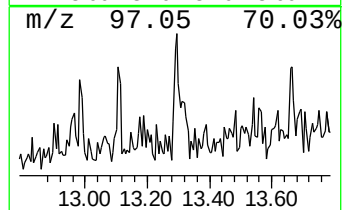
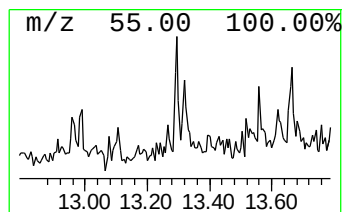
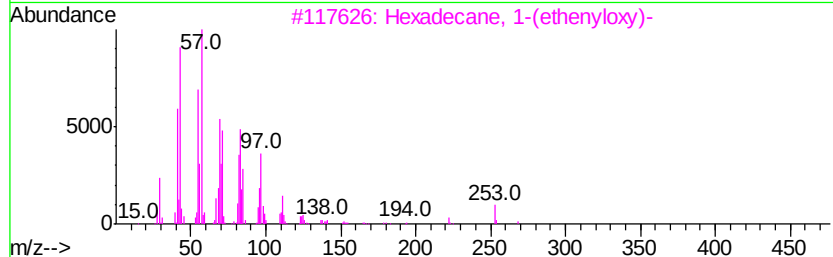
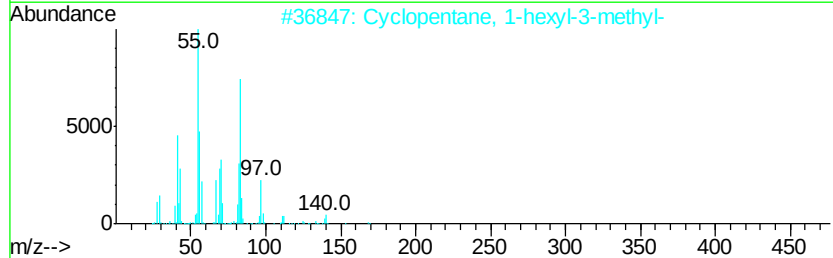
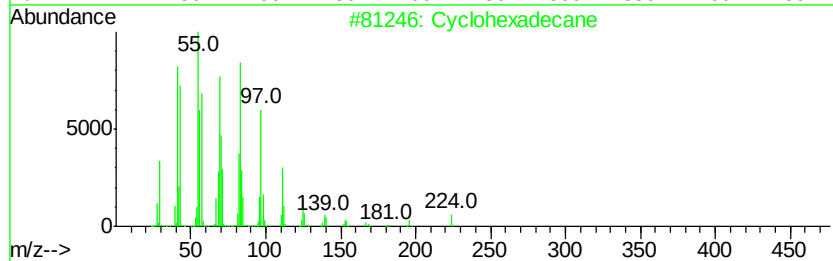
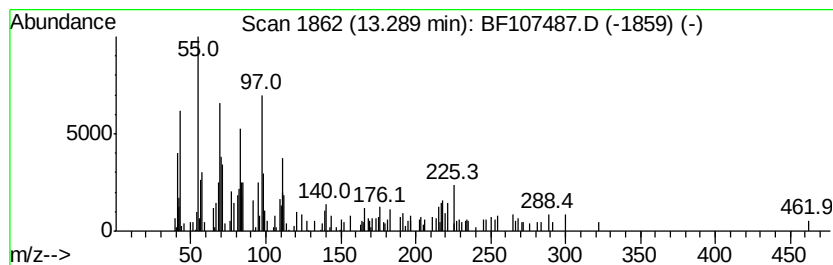
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclohexadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	2.87 ng	143519	Chrysene-d12	14.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	86
2		Cyclopentane, 1-hexyl-3-methyl-	168	C12H24	061142-68-5	64
3		Hexadecane, 1-(ethenvloxy)-	268	C18H36O	000822-28-6	64
4		1-Octadecanol	270	C18H38O	000112-92-5	58
5		Cyclotetradecane	196	C14H28	000295-17-0	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
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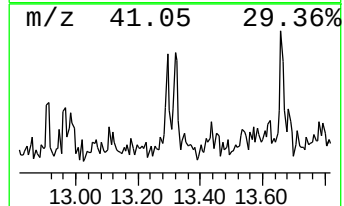
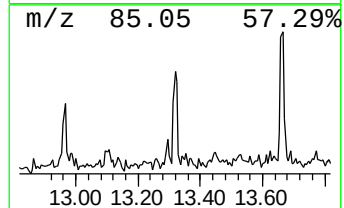
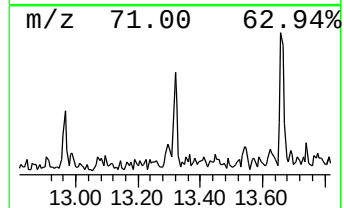
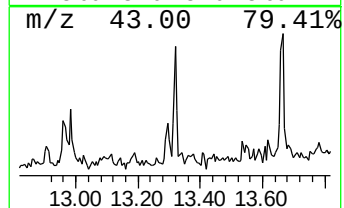
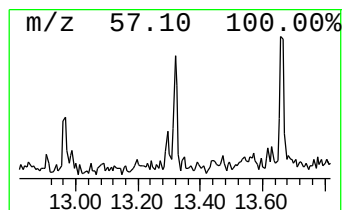
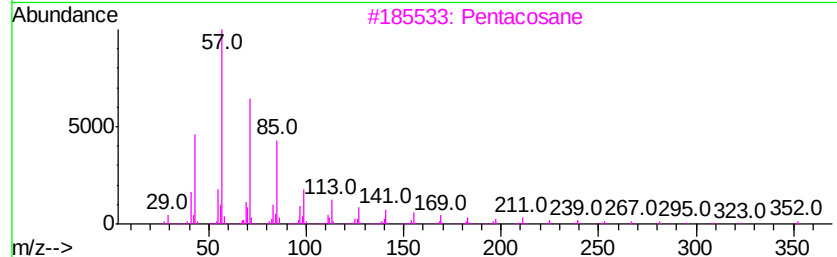
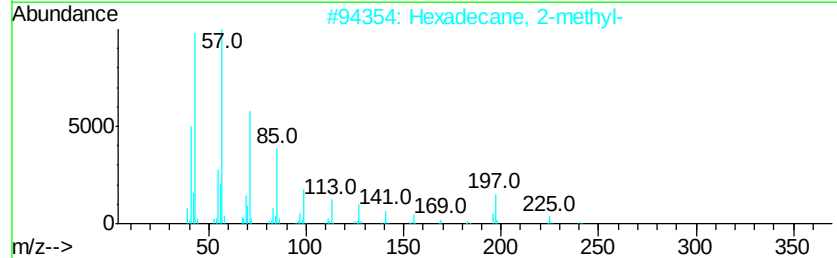
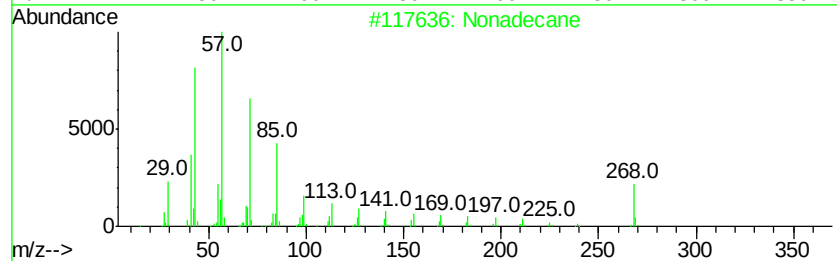
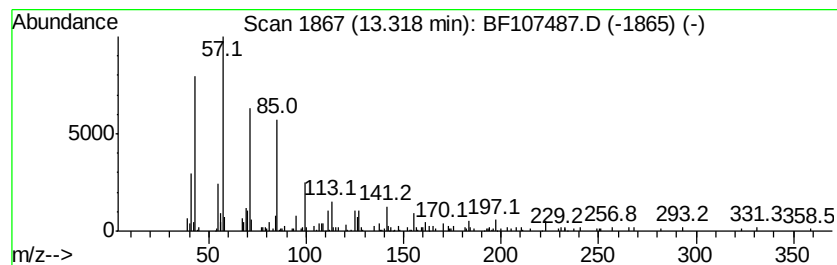
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	3.27 ng	163148	Chrysene-d12	14.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	91
3		Pentacosane	352	C25H52	000629-99-2	83
4		Heneicosane, 10-methyl-	310	C22H46	013287-25-7	80
5		Octacosane	394	C28H58	000630-02-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
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Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

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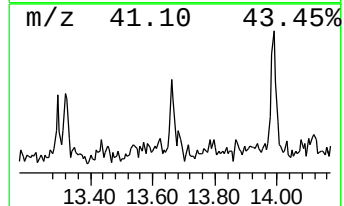
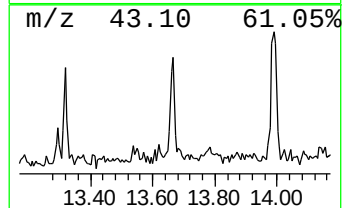
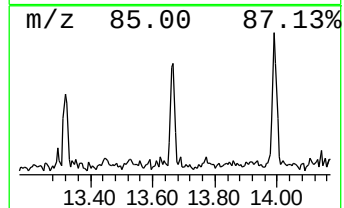
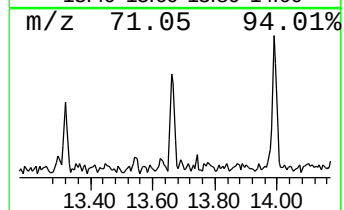
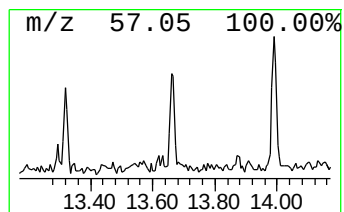
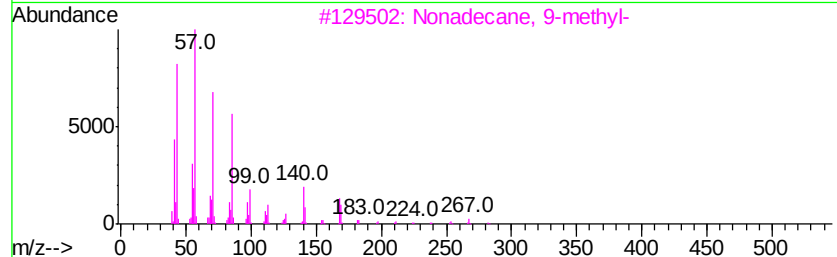
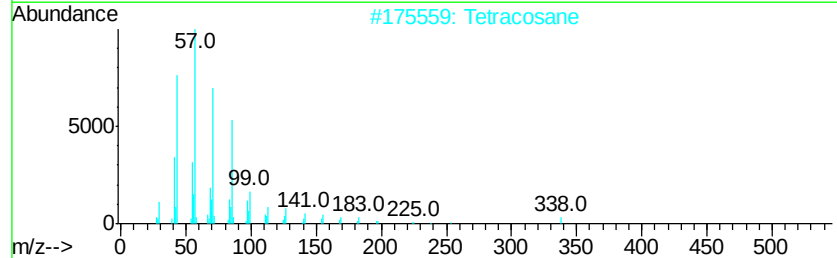
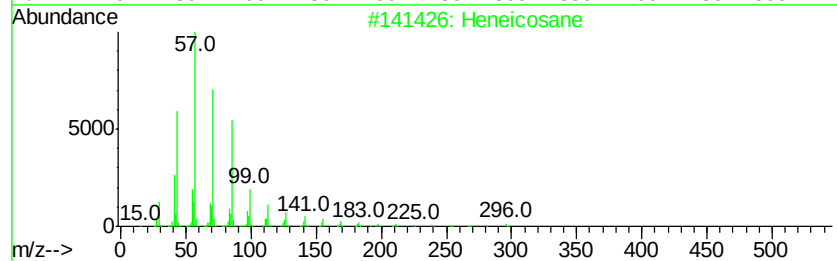
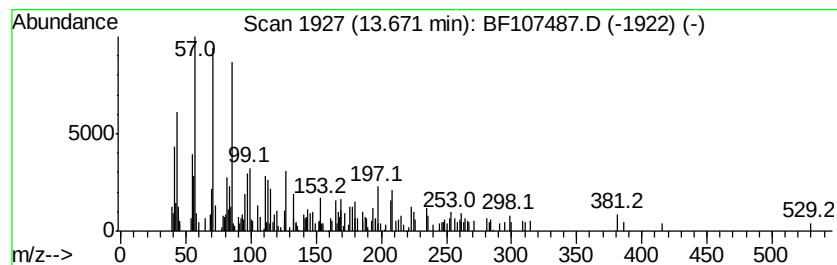
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	4.44 ng	221710	Chrysene-d12	14.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Tetracosane	338	C24H50	000646-31-1	86
3			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	83
4			Docosane	310	C22H46	000629-97-0	80
5			Octacosane	394	C28H58	000630-02-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
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Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

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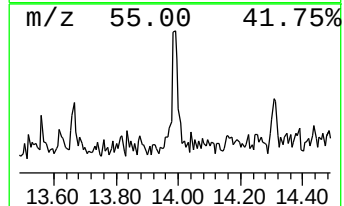
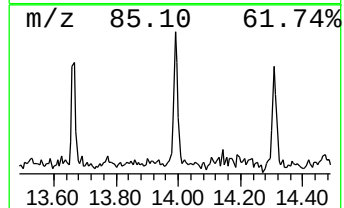
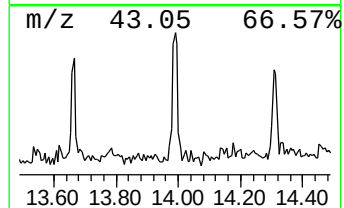
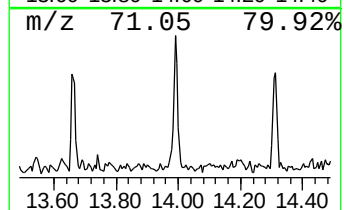
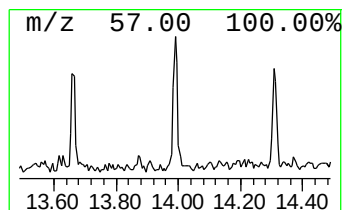
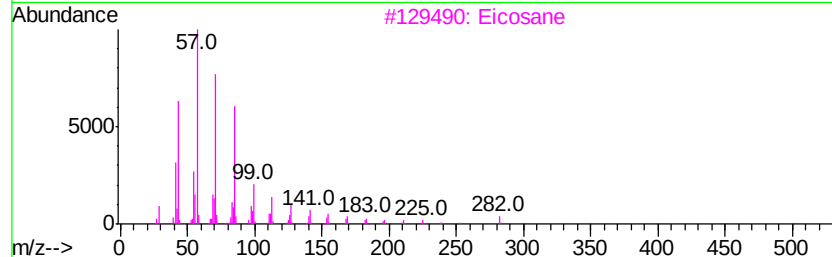
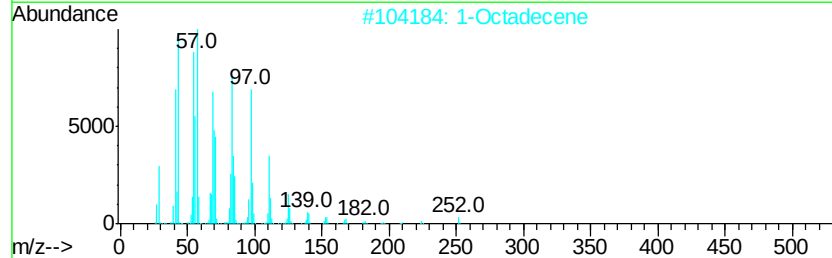
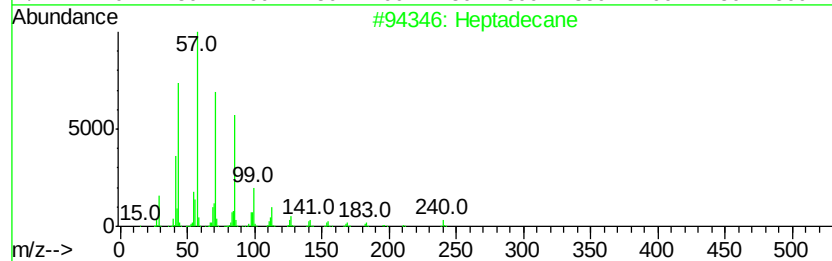
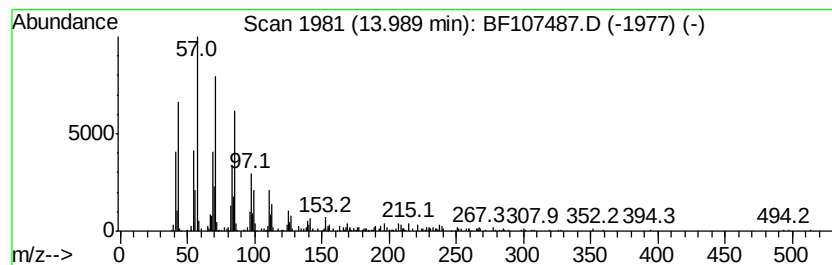
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	8.82 ng	440628	Chrysene-d12	14.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	93
2	1-Octadecene		252	C18H36	000112-88-9	92
3	Eicosane		282	C20H42	000112-95-8	89
4	Octacosane		394	C28H58	000630-02-4	86
5	Nonadecane, 9-methyl-		282	C20H42	013287-24-6	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
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Acq On : 18 Jul 2018 19:33
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Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

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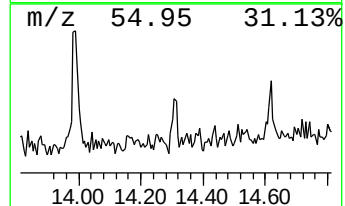
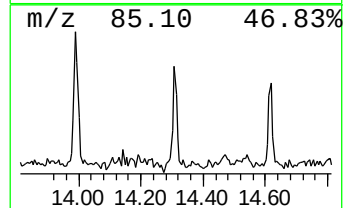
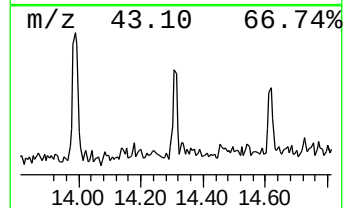
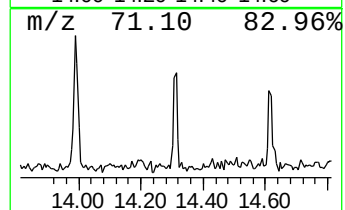
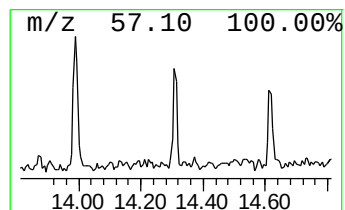
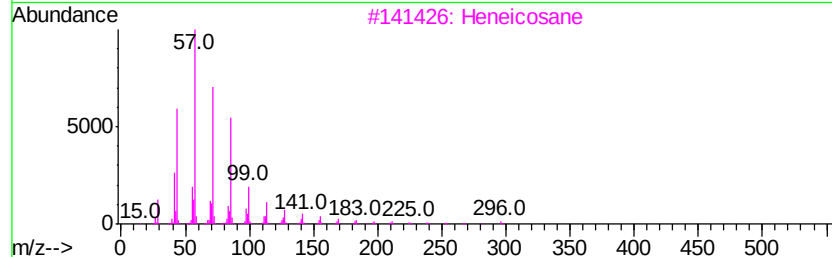
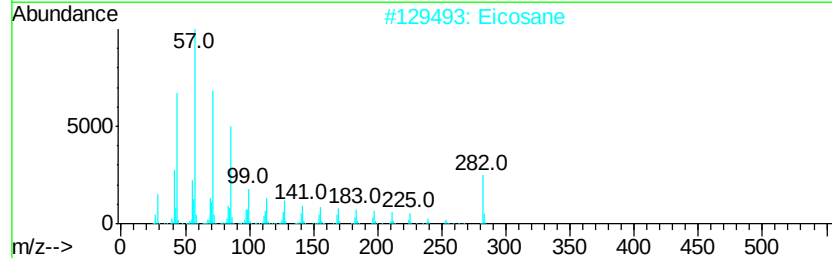
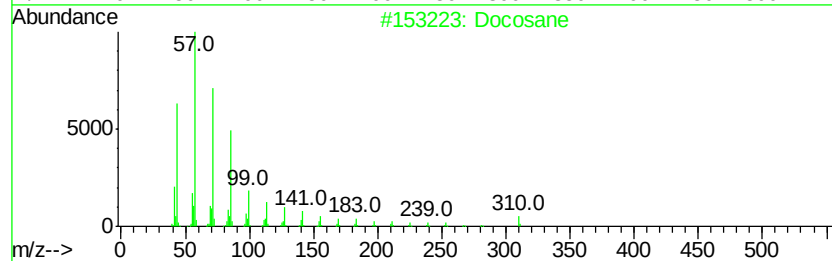
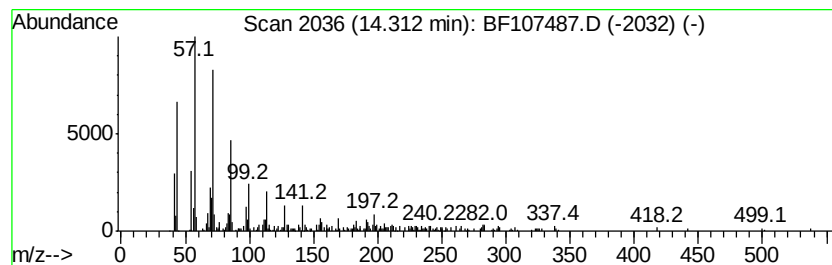
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.31	5.14 ng	256718	Chrysene-d12	14.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	90
2		Eicosane	282	C20H42	000112-95-8	90
3		Heneicosane	296	C21H44	000629-94-7	87
4		Triacontane	422	C30H62	000638-68-6	87
5		Pentacosane	352	C25H52	000629-99-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
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Sample : J4035-01
Misc :
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Instrument :
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ClientSampleId :
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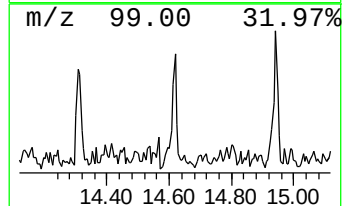
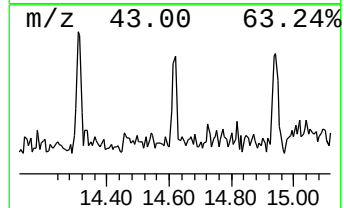
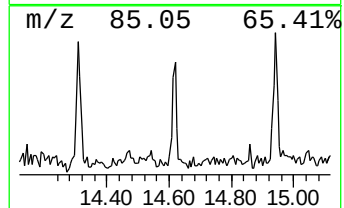
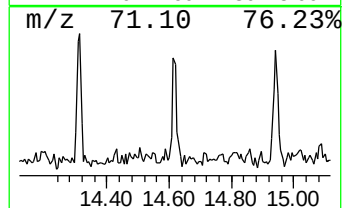
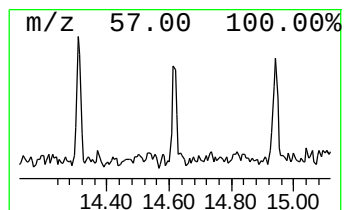
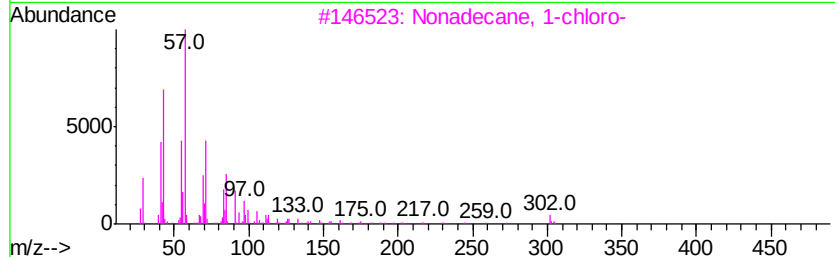
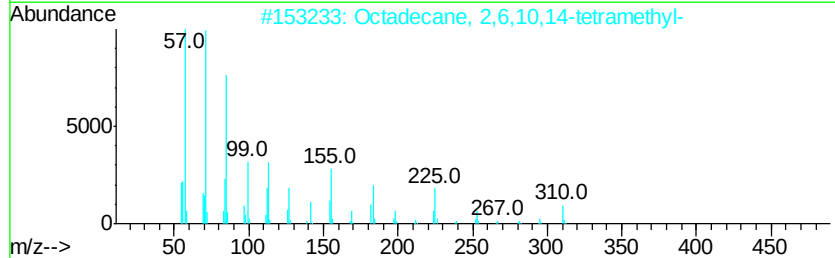
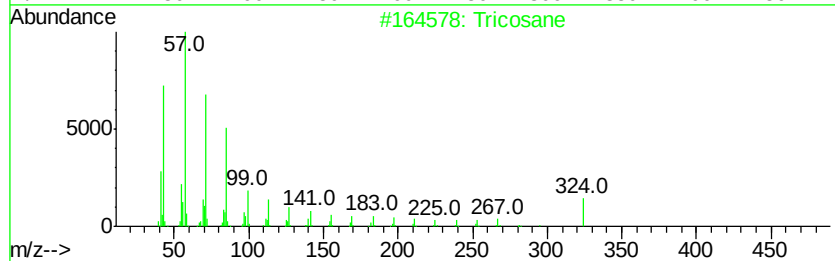
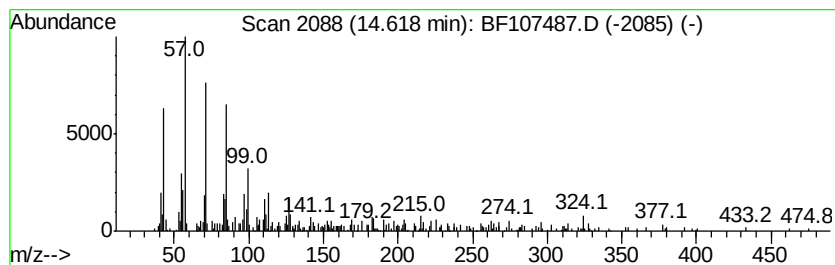
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Tricosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	4.25 ng	212241	Chrysene-d12	14.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	83
2		Octadecane, 2,6,10,14-tetramethyl-	310	C22H46	054964-82-8	76
3		Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	76
4		Docosane	310	C22H46	000629-97-0	74
5		Tetracontane	563	C40H82	004181-95-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
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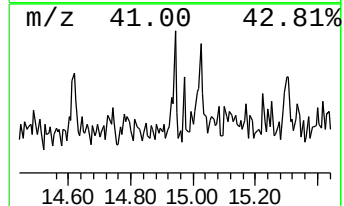
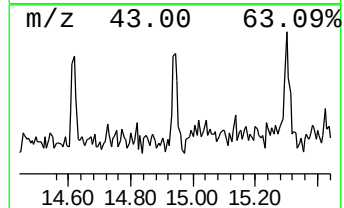
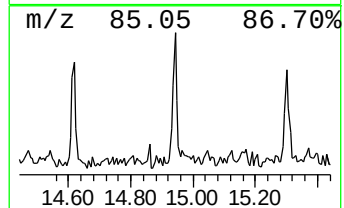
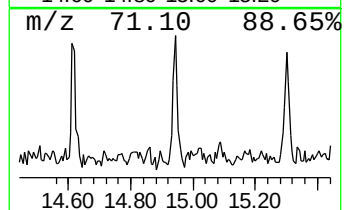
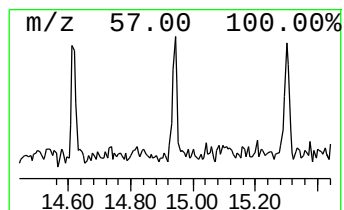
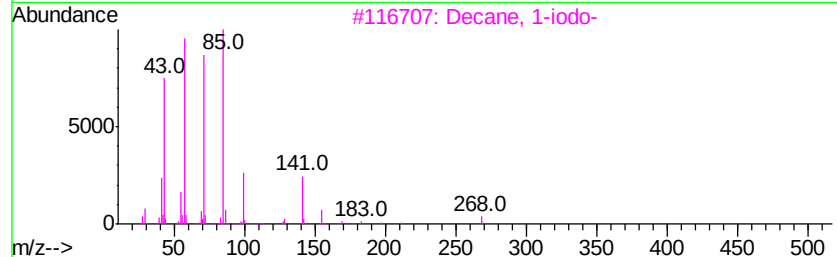
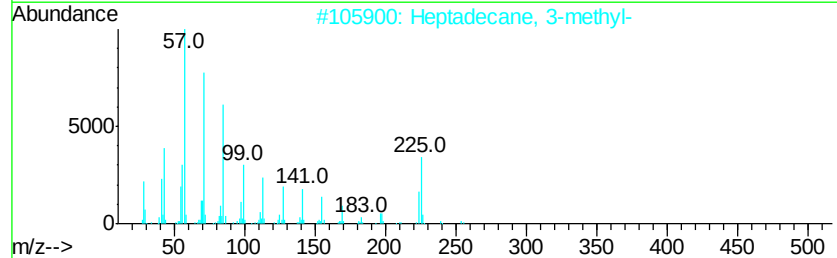
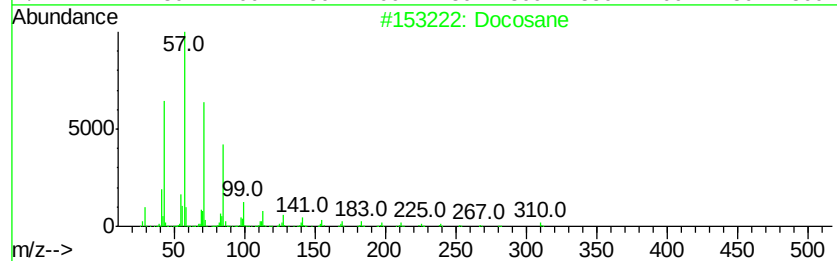
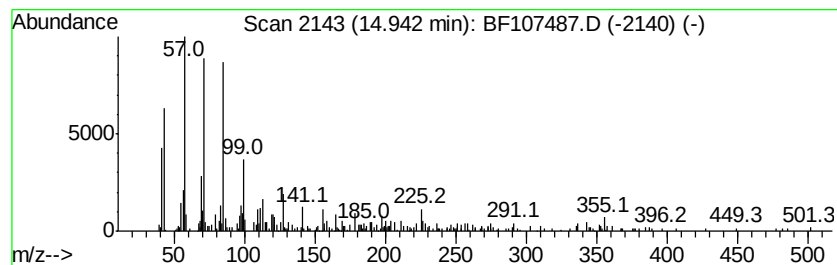
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Heptadecane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	3.75 ng	174128	Perylene-d12	15.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	93
2		Heptadecane, 3-methyl-	254	C18H38	006418-44-6	87
3		Decane, 1-iodo-	268	C10H21I	002050-77-3	87
4		Pentadecane	212	C15H32	000629-62-9	83
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
Operator : JU/SJ
Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2

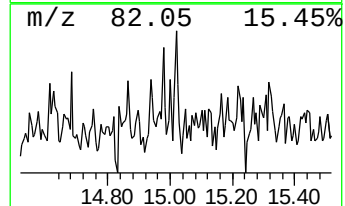
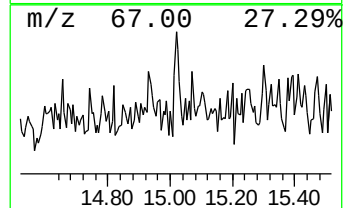
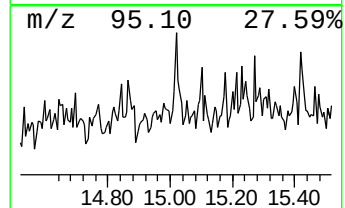
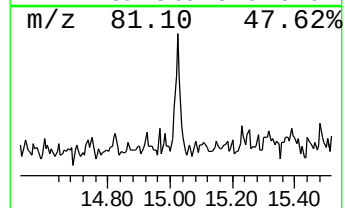
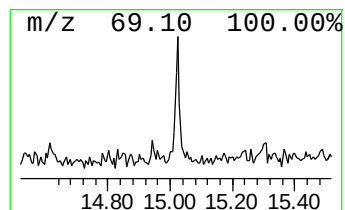
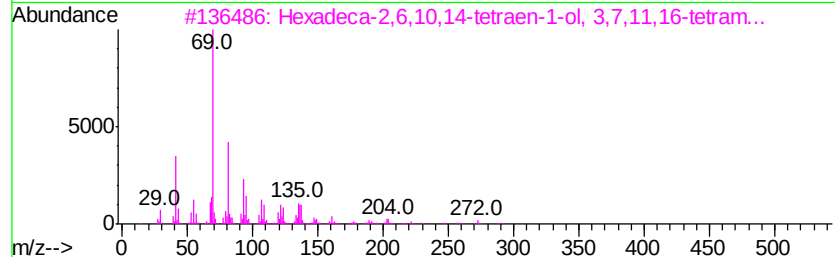
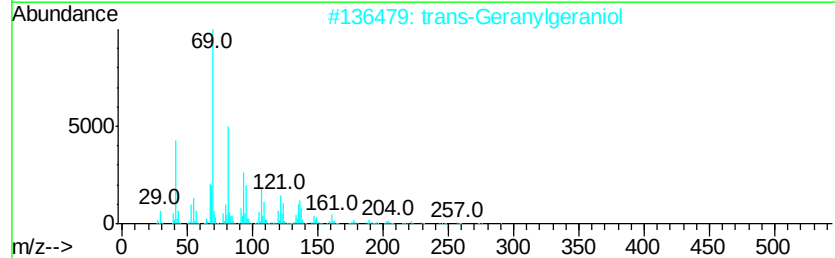
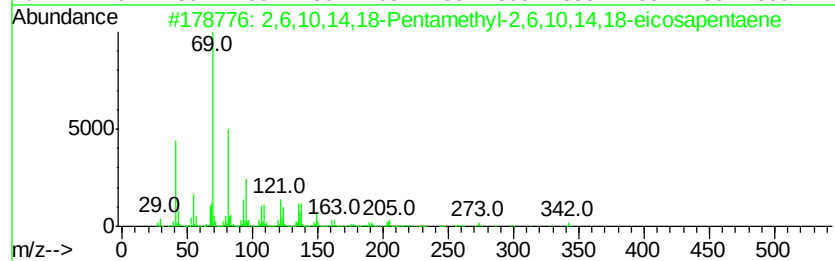
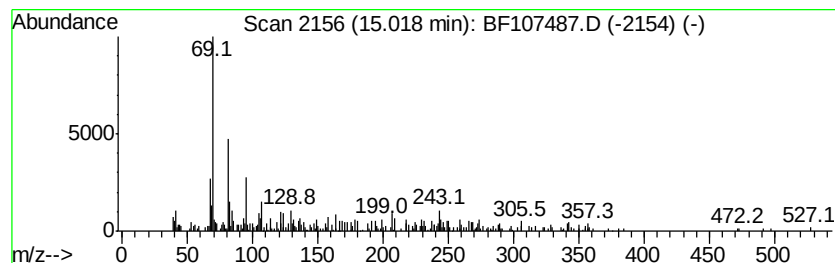
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2,6,10,14,18-Pentamethyl-2,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	2.91 ng	135465	Perylene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	81		
2	trans-Geranylgeraniol	290	C20H34O	024034-73-9	70		
3	Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	58		
4	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	53		
5	Trifluoroacetyl-lavandulol	250	C12H17F3O2	028673-24-7	53		



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
Operator : JU/SJ
Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2

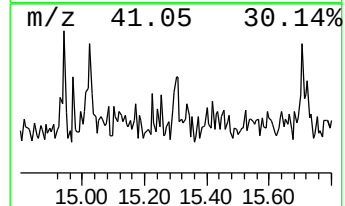
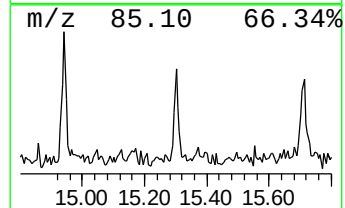
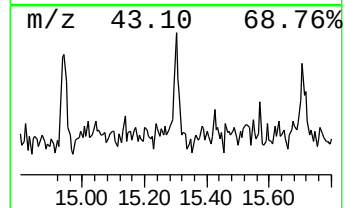
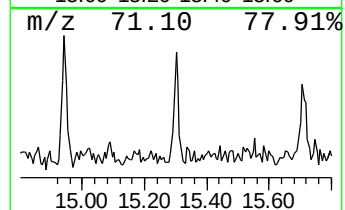
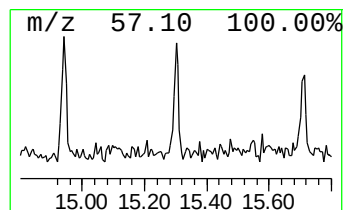
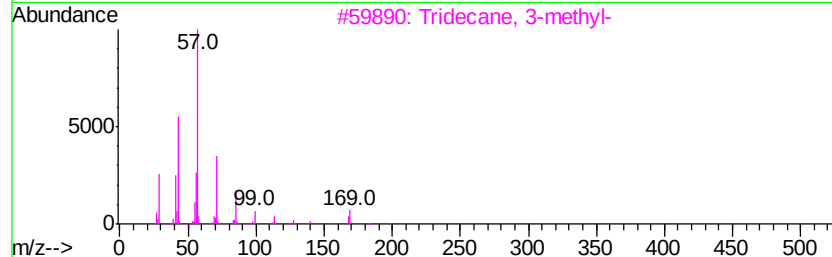
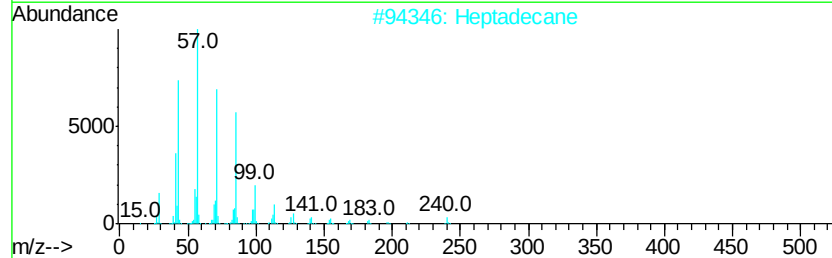
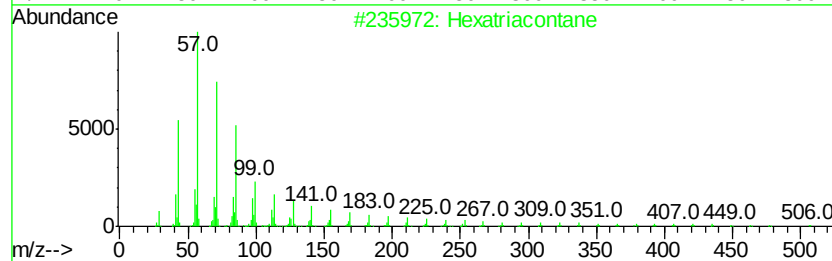
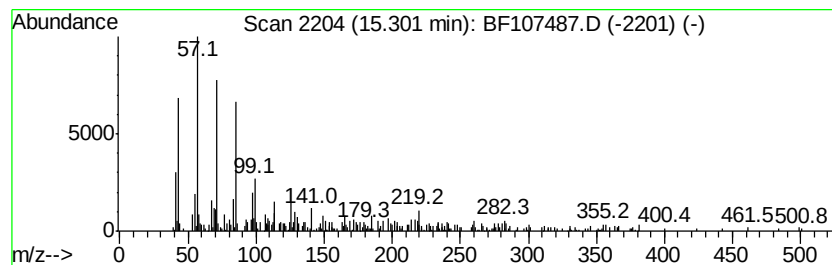
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexatriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.30	3.48 ng	161581	Perylene-d12	15.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexatriacontane	507	C36H74	000630-06-8	86
2		Heptadecane	240	C17H36	000629-78-7	81
3		Tridecane, 3-methyl-	198	C14H30	006418-41-3	80
4		Eicosane	282	C20H42	000112-95-8	74
5		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
Data File : BF107487.D
Acq On : 18 Jul 2018 19:33
Operator : JU/SJ
Sample : J4035-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2

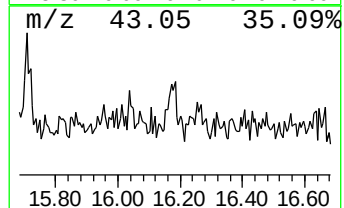
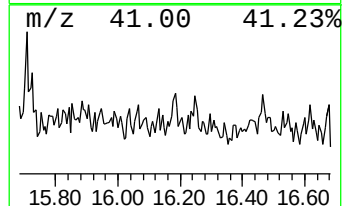
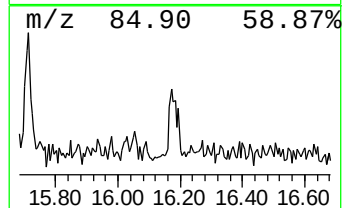
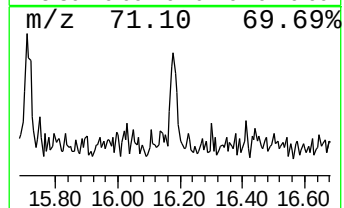
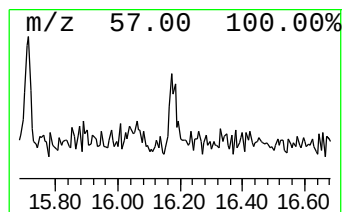
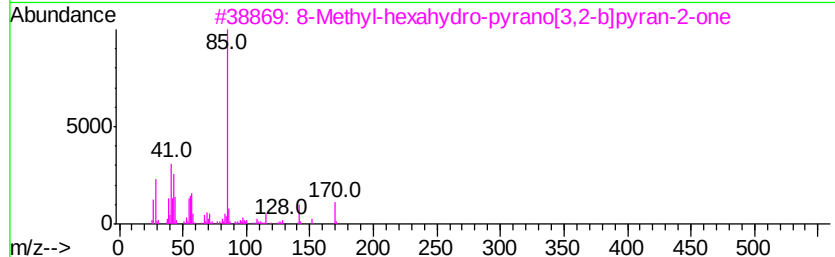
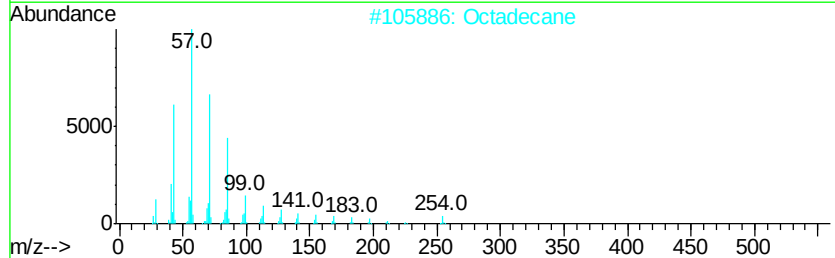
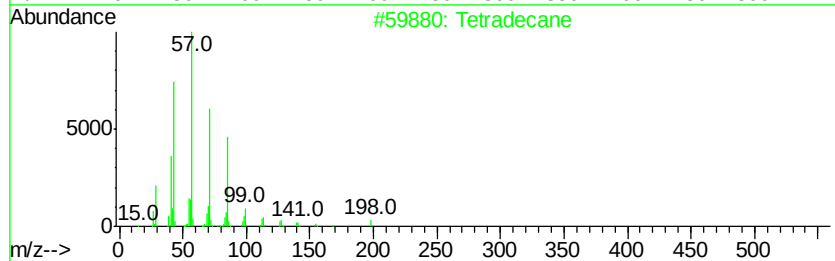
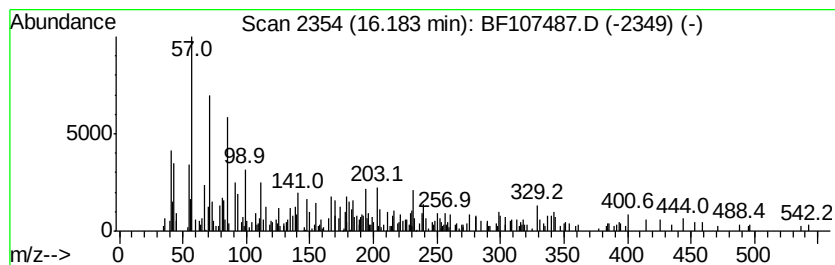
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.18 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	3.50 ng	162604	Perylene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	38
2			Octadecane	254	C18H38	000593-45-3	30
3			8-Methyl-hexahydro-pyrano[3,2-b]...	170	C9H14O3	1000190-63-9	27
4			Azacyclonol	267	C18H21NO	000115-46-8	27
5			Malonic acid, 3-methylpentyl oct...	300	C17H32O4	1000348-28-2	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071818\
 Data File : BF107487.D
 Acq On : 18 Jul 2018 19:33
 Operator : JU/SJ
 Sample : J4035-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071618.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-Pentanone, 4-hy...	5.23	8.5	ng	466360	1	6.98	1101000	20.0
unknown6.76	6.76	71.3	ng	3923230	1	6.98	1101000	20.0
Pentaethylene glycol	9.97	3.4	ng	241251	3	10.03	1431450	20.0
n-Nonadecanol-1	12.55	5.1	ng	330933	4	11.52	1289200	20.0
Octadecane, 1-chl...	12.97	2.3	ng	115889	5	14.17	998794	20.0
unknown12.98	12.98	3.0	ng	151386	5	14.17	998794	20.0
Cyclohexadecane	13.29	2.9	ng	143519	5	14.17	998794	20.0
Nonadecane	13.32	3.3	ng	163148	5	14.17	998794	20.0
Heneicosane	13.67	4.4	ng	221710	5	14.17	998794	20.0
Heptadecane	13.99	8.8	ng	440628	5	14.17	998794	20.0
Docosane	14.31	5.1	ng	256718	5	14.17	998794	20.0
Tricosane	14.62	4.3	ng	212241	5	14.17	998794	20.0
Heptadecane, 3-me...	14.94	3.8	ng	174128	6	15.71	929871	20.0
2,6,10,14,18-Pent...	15.02	2.9	ng	135465	6	15.71	929871	20.0
Hexatriacontane	15.30	3.5	ng	161581	6	15.71	929871	20.0
unknown16.18	16.18	3.5	ng	162604	6	15.71	929871	20.0