

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
 Data File : BF109911.D
 Acq On : 16 Oct 2018 17:53
 Operator : JU/SJ
 Sample : J5516-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1012-S-DH-29

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-BF101518.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	2.116	3	5	11	rVB	373984	416261	13.08%	1.260%
2	2.322	33	40	46	rVB	557933	746390	23.45%	2.260%
3	5.210	527	531	538	rVB	113807	135874	4.27%	0.411%
4	5.598	592	597	600	rBV	3627008	3126615	98.23%	9.467%
5	6.592	761	766	769	rBV	3581030	3094390	97.22%	9.370%
6	6.745	788	792	795	rBV	3775070	3182977	100.00%	9.638%
7	6.981	828	832	835	rBV	1635038	1227626	38.57%	3.717%
8	7.133	855	858	862	rVB	2867808	2591124	81.41%	7.846%
9	7.539	922	927	930	rBV	2310946	1901957	59.75%	5.759%
10	8.263	1046	1050	1053	rBV	1864132	1454063	45.68%	4.403%
11	9.333	1228	1232	1235	rBV	3758735	3088563	97.03%	9.352%
12	9.727	1295	1299	1302	rBV	188297	167471	5.26%	0.507%
13	10.021	1345	1349	1353	rBV	1527046	1295156	40.69%	3.922%
14	10.810	1479	1483	1486	rBV	1617371	1369162	43.02%	4.146%
15	11.510	1599	1602	1605	rBV	1281833	1086749	34.14%	3.291%
16	12.904	1837	1839	1843	rVB	45319	37272	1.17%	0.113%
17	13.098	1867	1872	1875	rBV	2912182	2656516	83.46%	8.044%
18	13.568	1947	1952	1956	rBV	366713	528566	16.61%	1.600%
19	13.598	1956	1957	1962	rVB	149910	157459	4.95%	0.477%
20	13.974	2018	2021	2027	rBV2	181392	213352	6.70%	0.646%
21	14.157	2048	2052	2055	rBV	1705406	1496393	47.01%	4.531%
22	14.251	2065	2068	2074	rBV4	50589	77920	2.45%	0.236%
23	14.298	2074	2076	2082	rVB	93979	86462	2.72%	0.262%
24	14.604	2125	2128	2133	rVB	153976	153679	4.83%	0.465%
25	14.927	2179	2183	2189	rVB	207284	217655	6.84%	0.659%
26	15.009	2194	2197	2202	rBV	83764	90461	2.84%	0.274%
27	15.286	2240	2244	2252	rVB	201539	234794	7.38%	0.711%
28	15.686	2306	2312	2322	rVB	1311063	1769648	55.60%	5.358%
29	16.156	2387	2392	2396	rBV	155362	229029	7.20%	0.693%
30	16.698	2479	2484	2492	rVB2	106154	192229	6.04%	0.582%

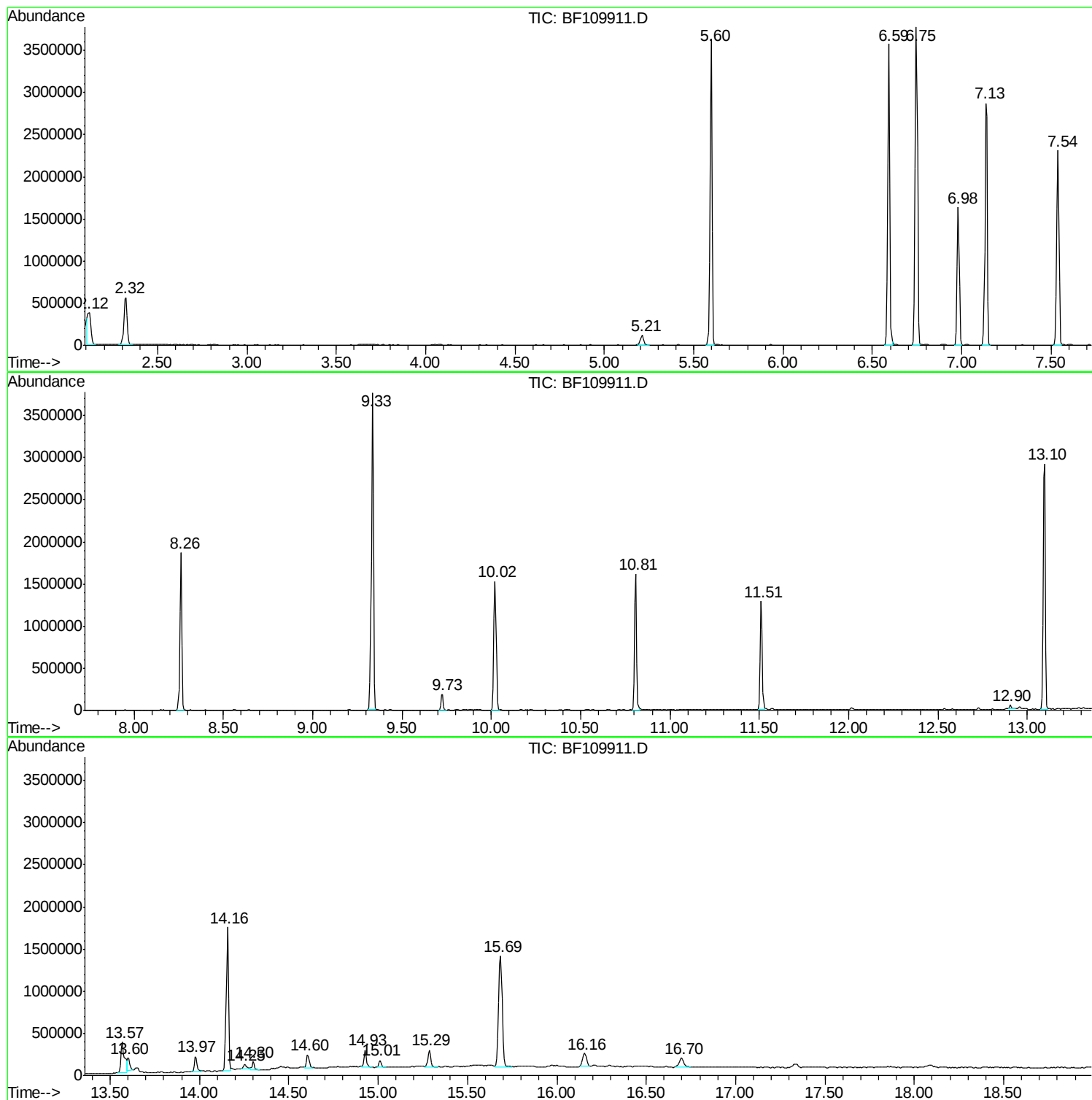
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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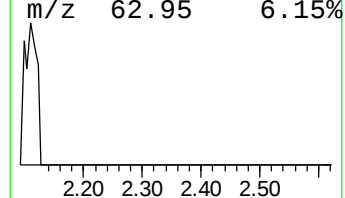
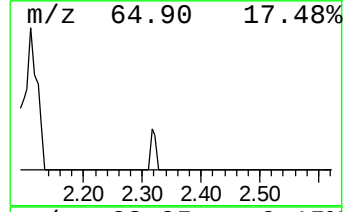
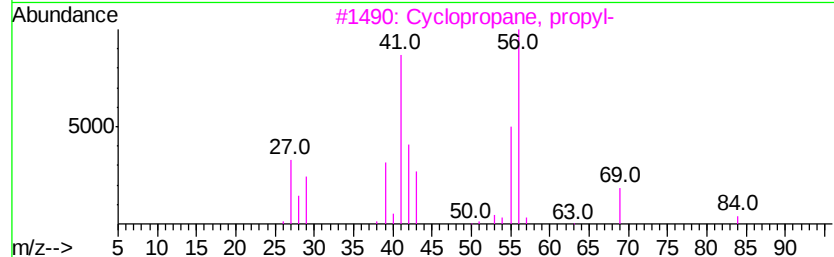
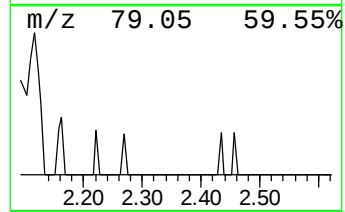
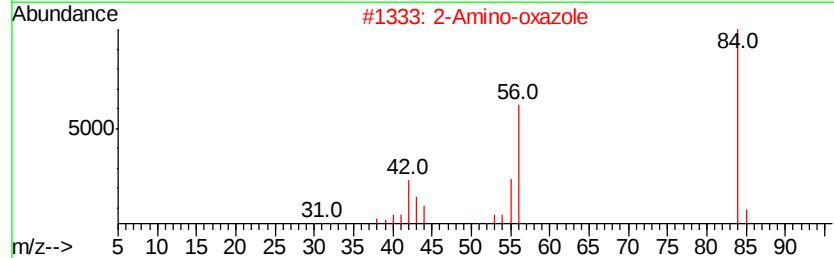
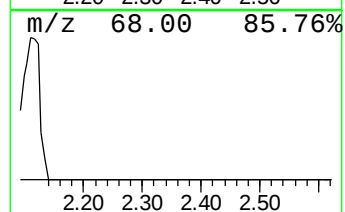
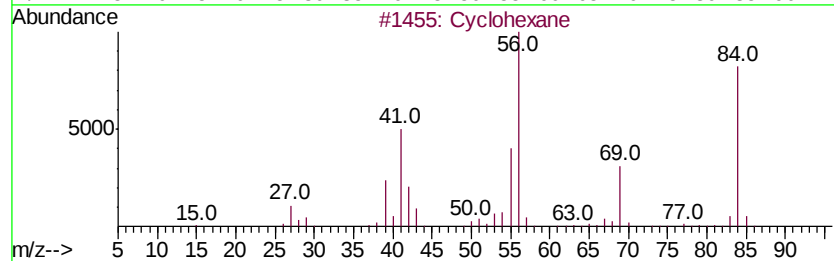
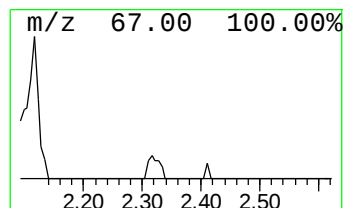
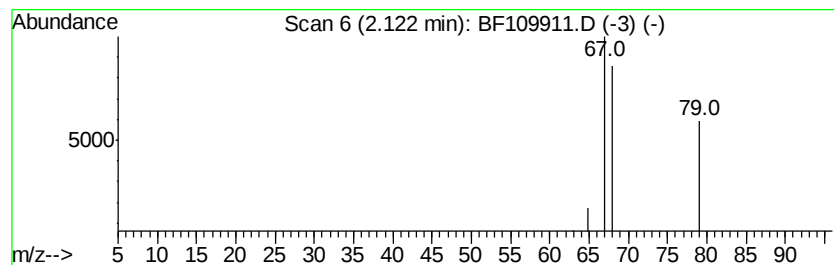
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	6.78 ng	416261	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	87
2		2-Amino-oxazole	84	C3H4N2O	004570-45-0	50
3		Cyclopropane, propyl-	84	C6H12	002415-72-7	37
4		1-Hexene, 5-methyl-	98	C7H14	003524-73-0	25
5		Oxirane, (1-methylbutyl)-	114	C7H14O	053229-39-3	25



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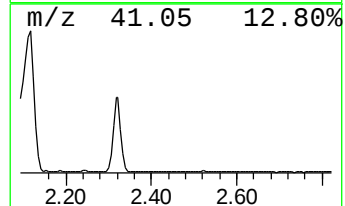
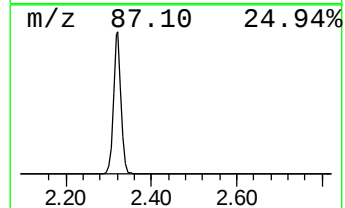
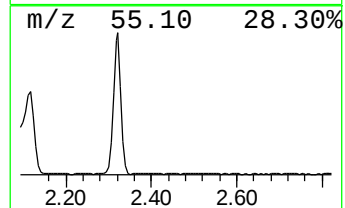
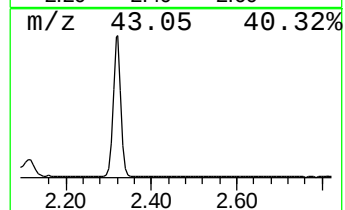
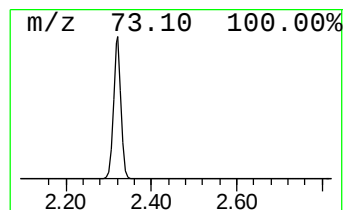
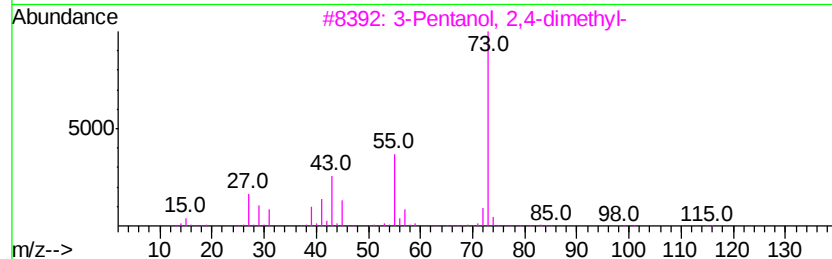
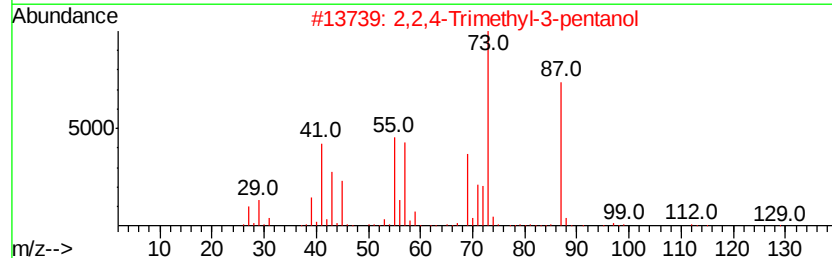
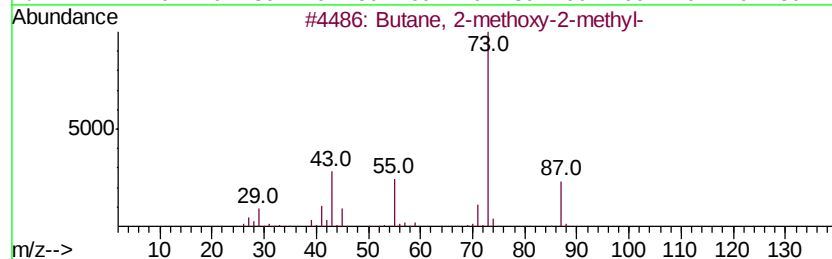
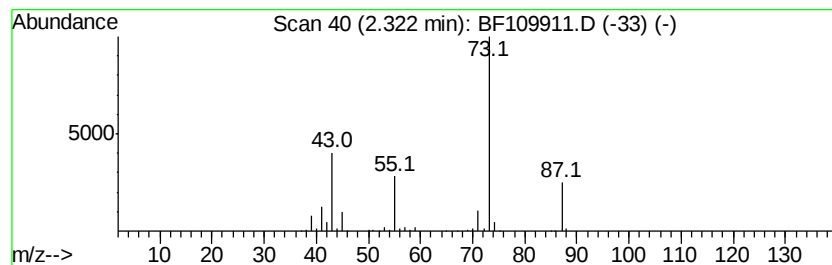
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	12.16 ng	746390	1,4-Dichlorobenzene-d4	6.98

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		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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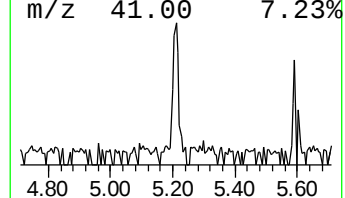
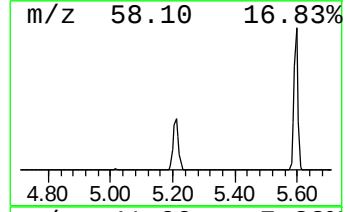
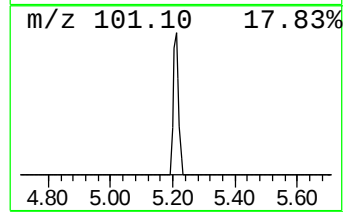
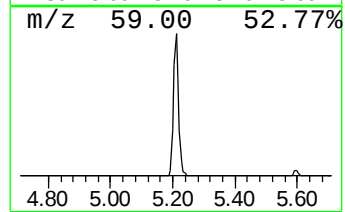
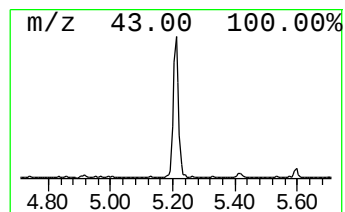
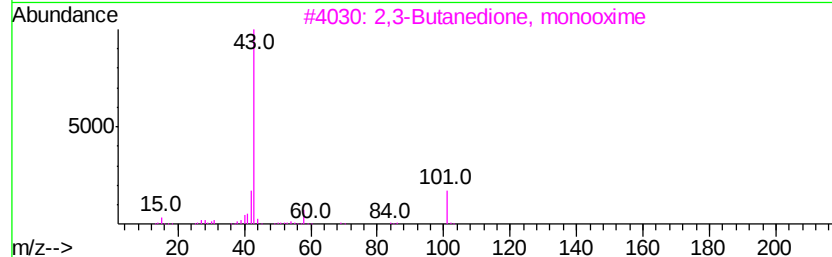
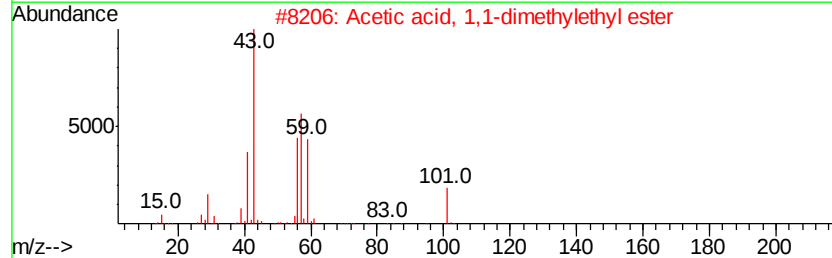
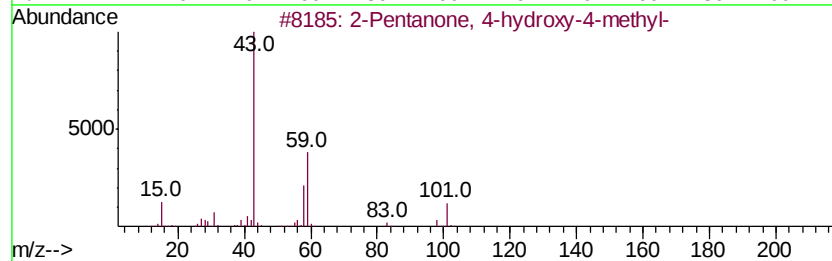
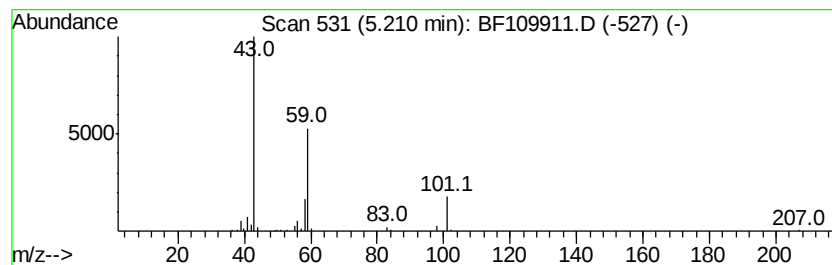
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	2.21 ng	135874	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetone	58	C3H6O	000067-64-1	9



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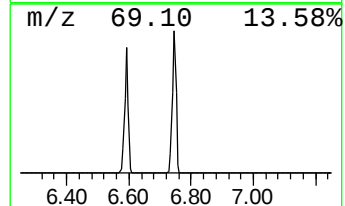
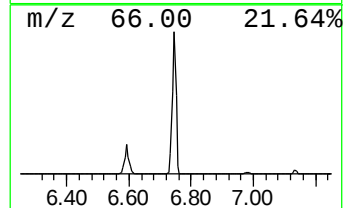
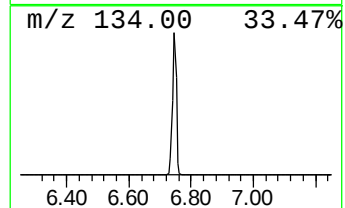
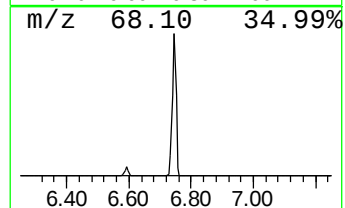
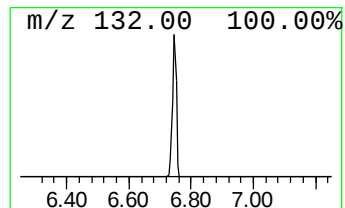
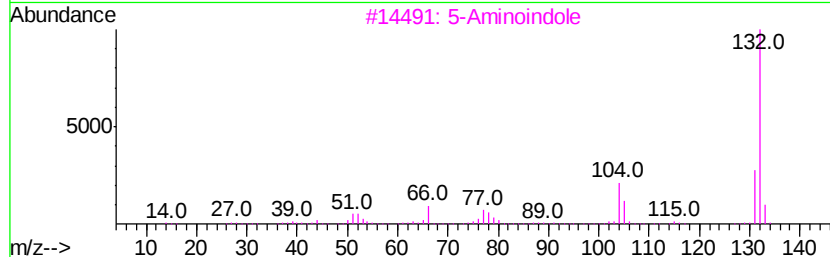
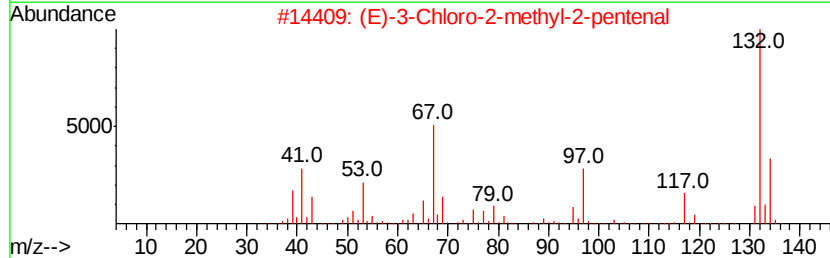
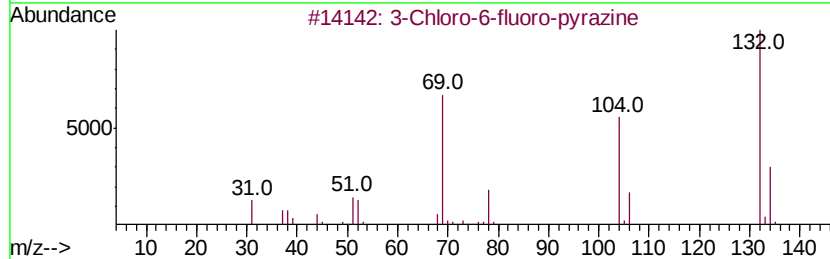
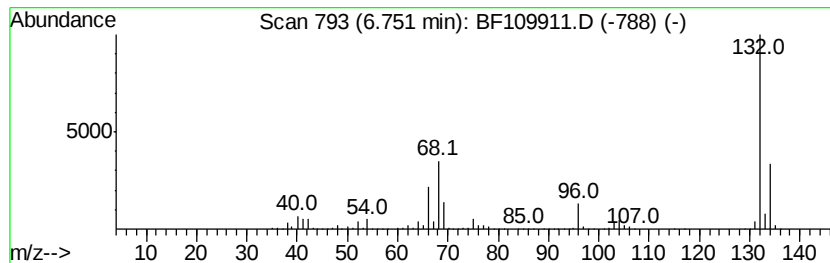
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Peak Number 4 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	51.86 ng	3182980	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3			5-Aminoindole	132	C8H8N2	005192-03-0	12
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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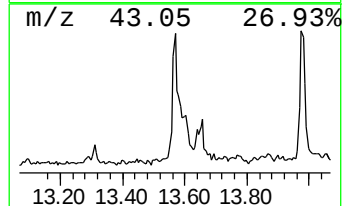
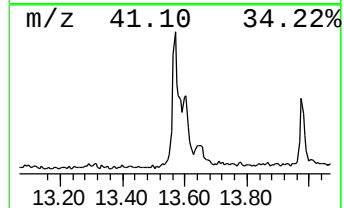
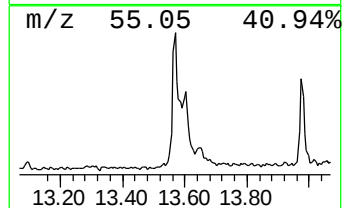
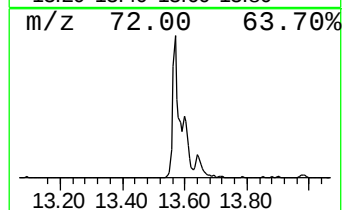
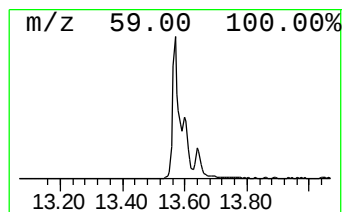
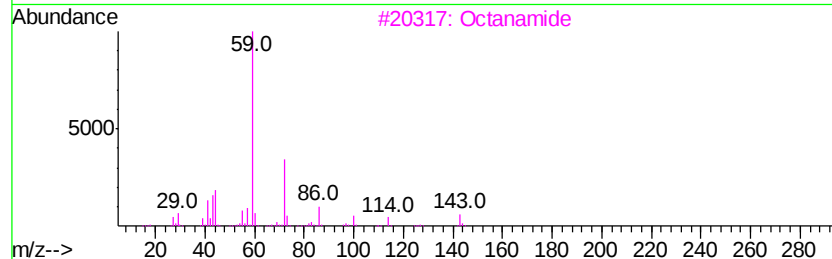
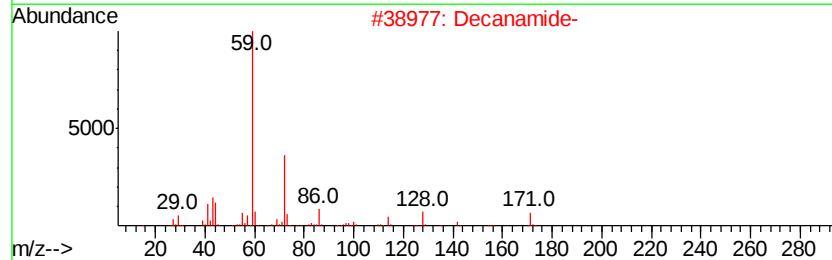
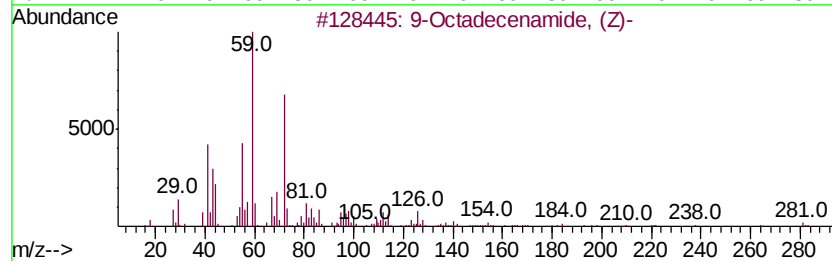
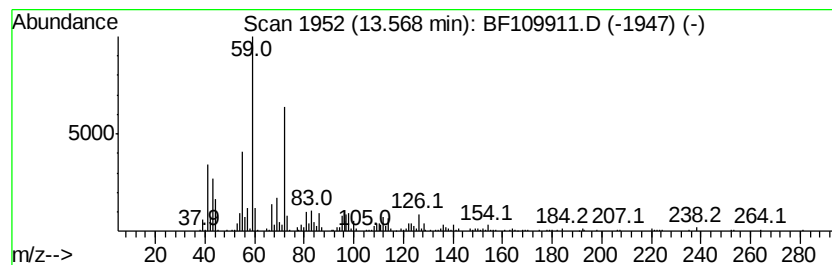
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Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	7.06 ng	528566	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		Decanamide-	171	C10H21NO	002319-29-1	64
3		Octanamide	143	C8H17NO	000629-01-6	59
4		Hexadecanamide	255	C16H33NO	000629-54-9	56
5		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	50



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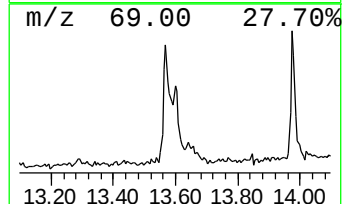
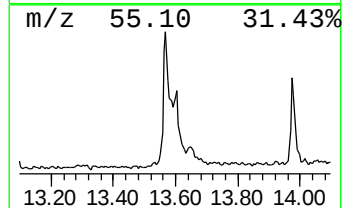
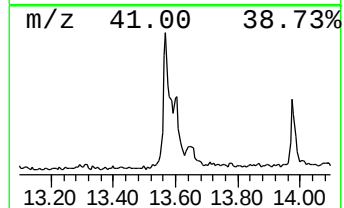
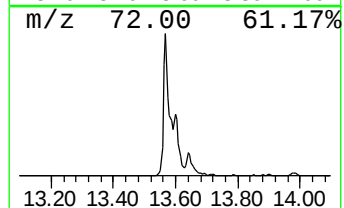
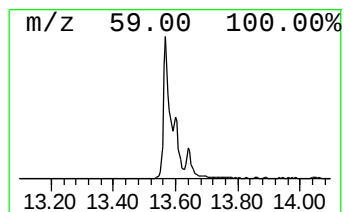
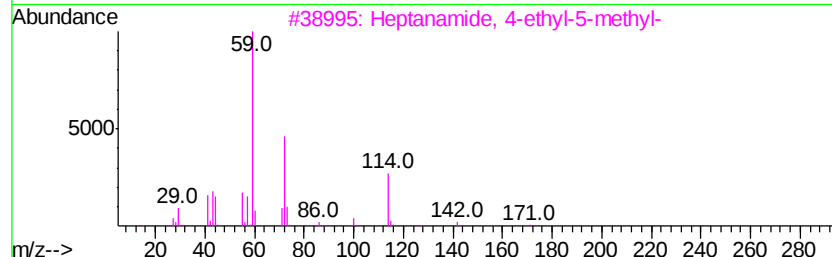
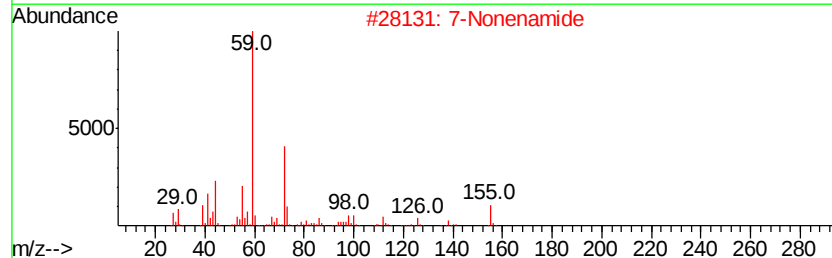
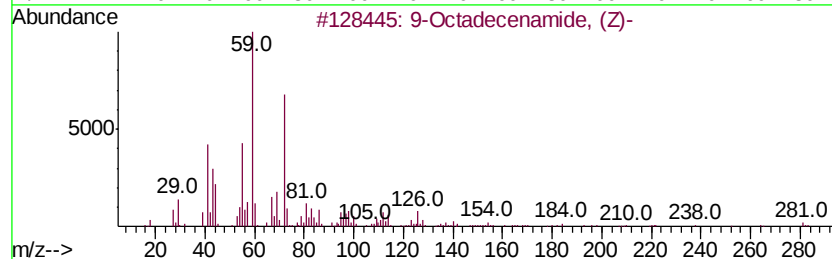
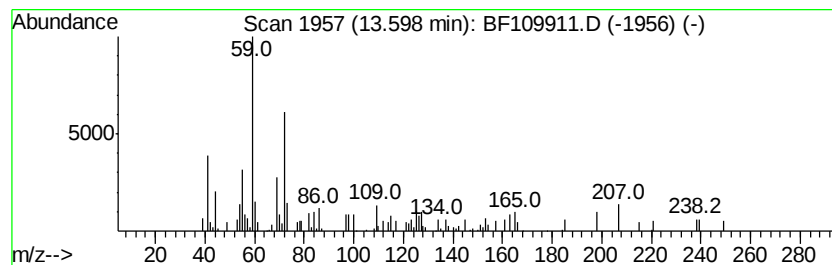
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 7-Nonenamide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	2.10 ng	157459	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	52
2		7-Nonenamide	155	C9H17NO	090949-53-4	43
3		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	40
4		Nonanamide	157	C9H19NO	001120-07-6	37
5		Dodecanamide	199	C12H25NO	001120-16-7	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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Acq On : 16 Oct 2018 17:53
Operator : JU/SJ
Sample : J5516-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSSI-1012-S-DH-29

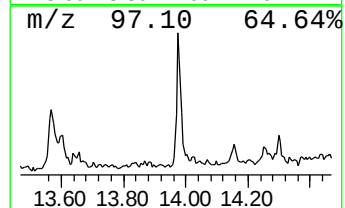
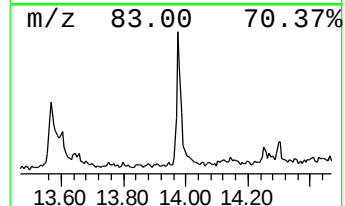
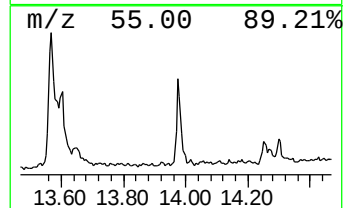
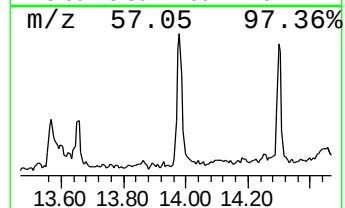
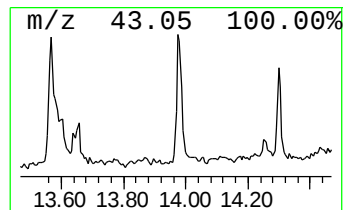
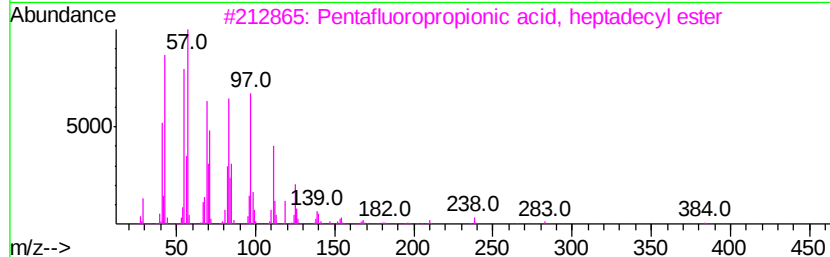
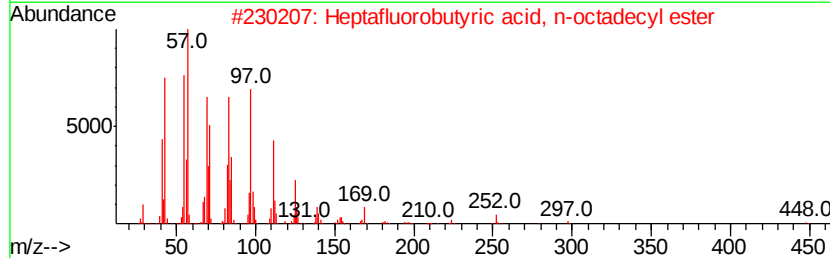
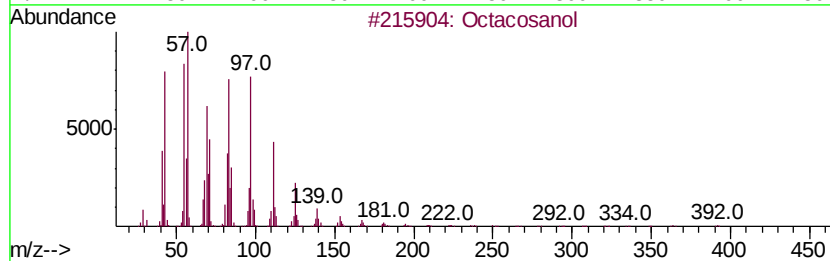
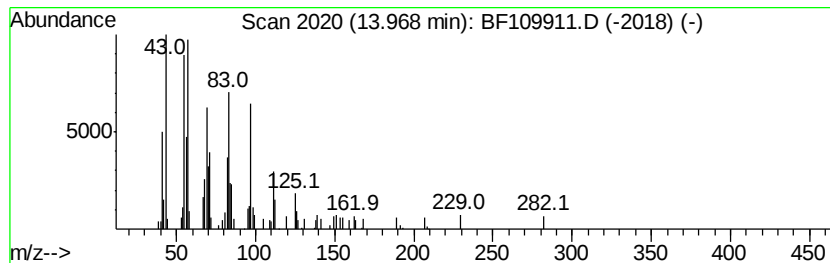
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	2.85 ng	213352	Chrysene-d12	14.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosanol	410	C28H58O	000557-61-9	94
2		Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	94
3		Pentafluoropropionic acid, heptadecyl ester	402	C20H35F5O2	959218-78-1	94
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	94



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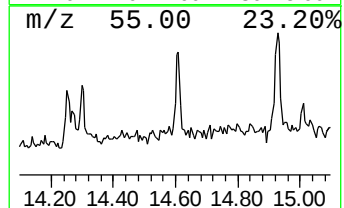
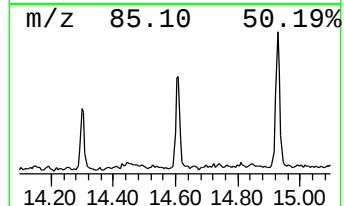
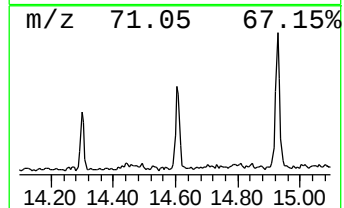
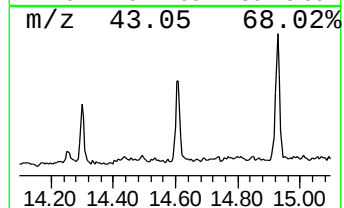
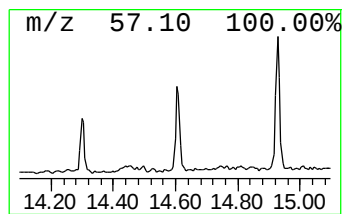
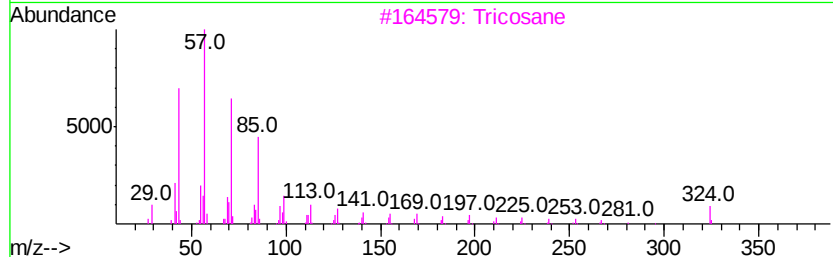
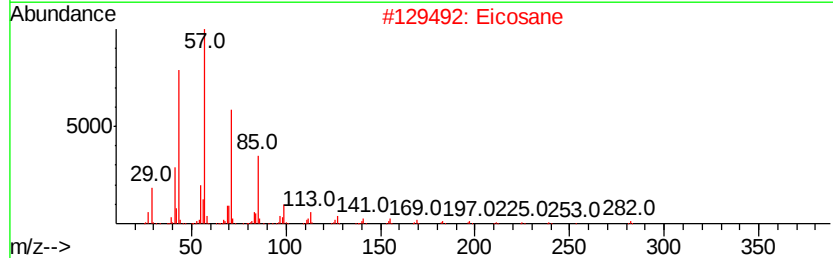
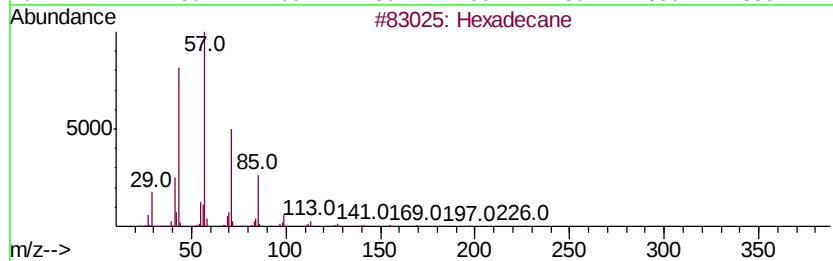
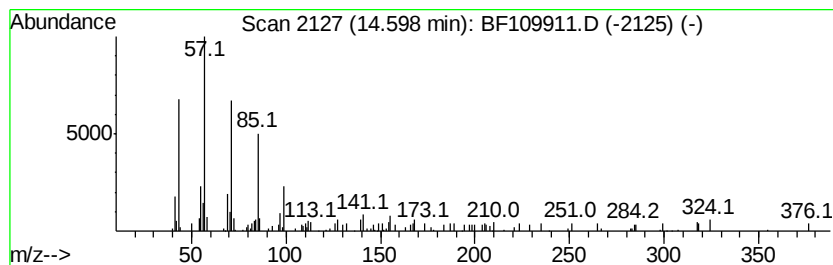
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.05 ng	153679	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	95
2		Eicosane	282	C20H42	000112-95-8	94
3		Tricosane	324	C23H48	000638-67-5	94
4		Tetratetracontane	619	C44H90	007098-22-8	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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Sample : J5516-08
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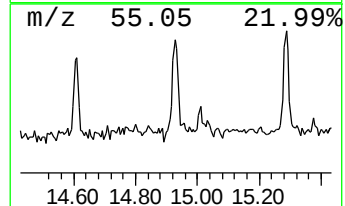
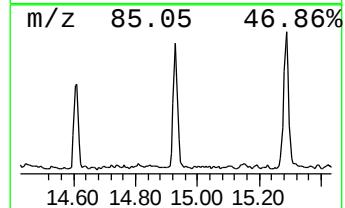
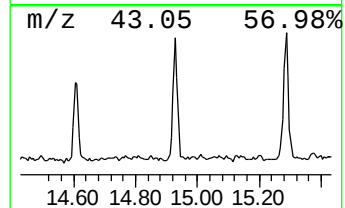
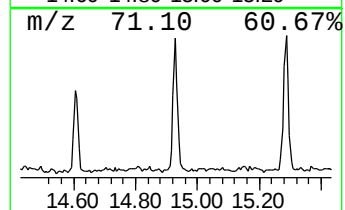
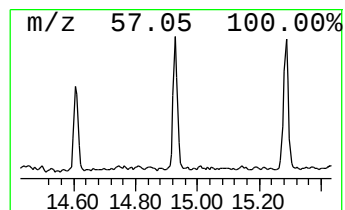
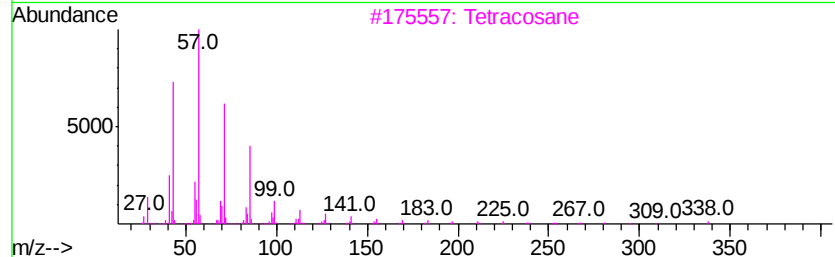
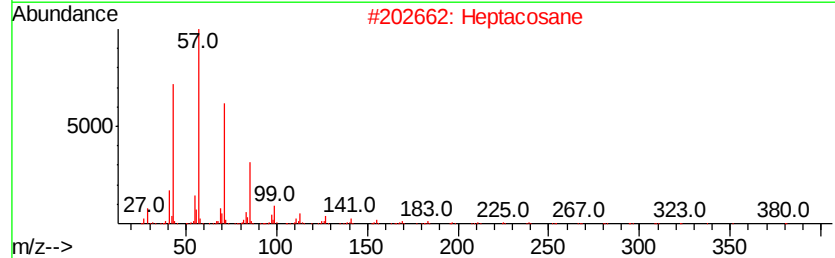
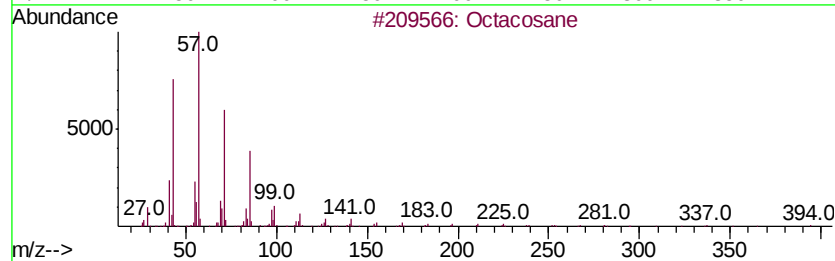
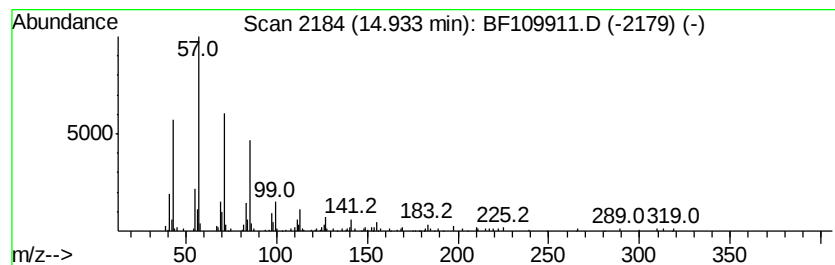
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	2.46 ng	217655	Perylene-d12	15.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	91
2	Heptacosane	380 C27H56	000593-49-7	91
3	Tetracosane	338 C24H50	000646-31-1	91
4	Pentadecane, 8-hexyl-	296 C21H44	013475-75-7	90
5	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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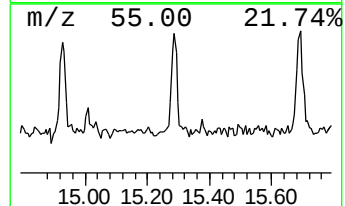
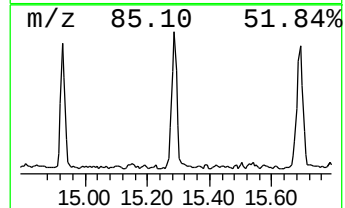
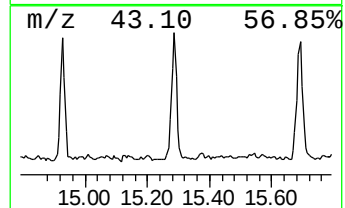
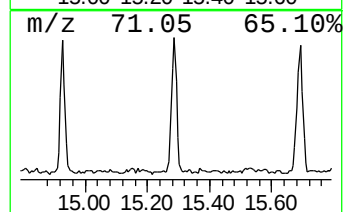
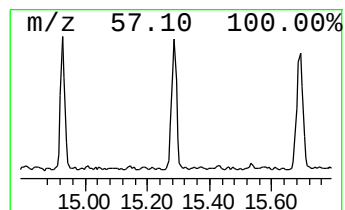
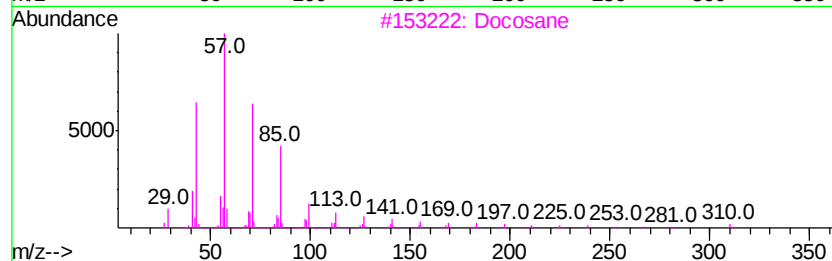
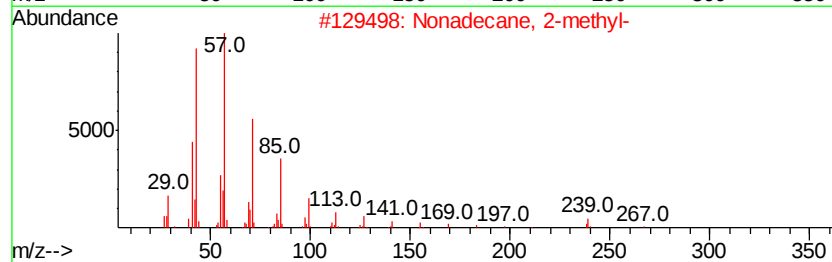
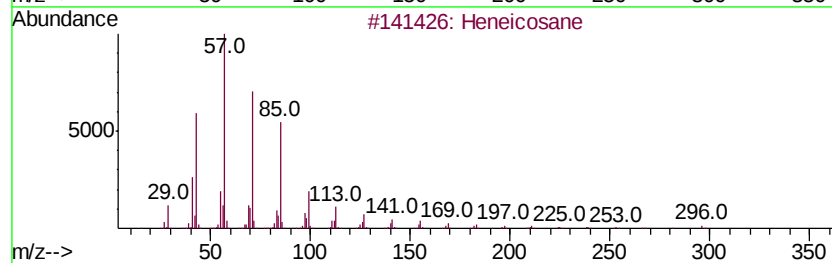
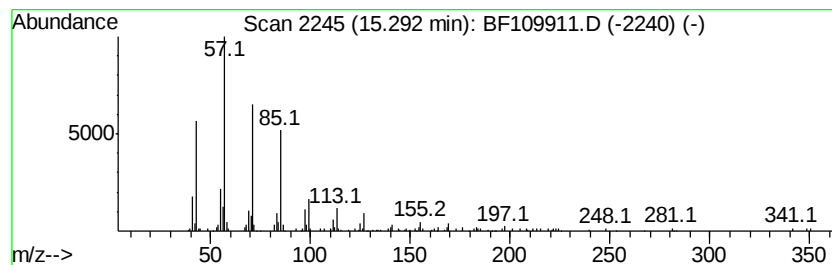
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	2.65 ng	234794	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	90
3		Docosane	310	C22H46	000629-97-0	90
4		Triacontane	422	C30H62	000638-68-6	90
5		Tetratetracontane	619	C44H90	007098-22-8	87



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Operator : JU/SJ
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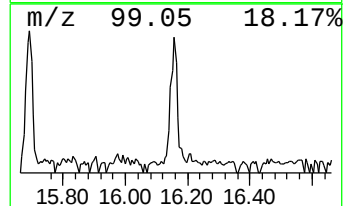
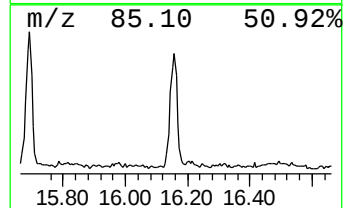
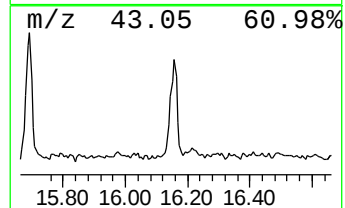
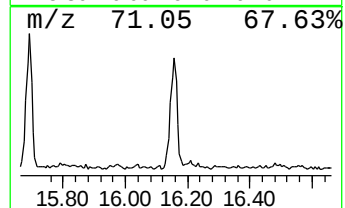
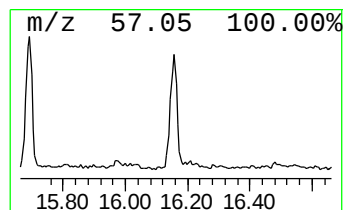
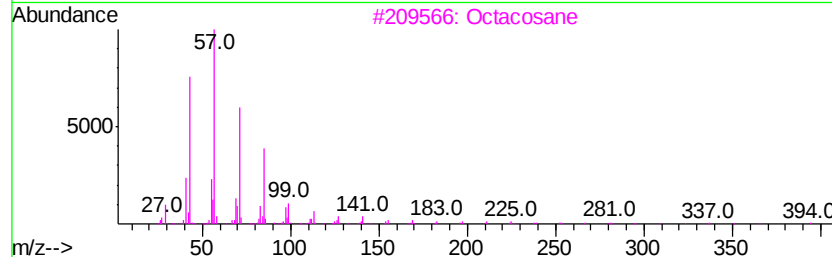
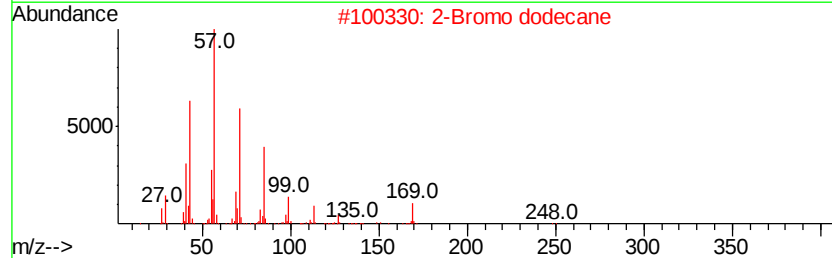
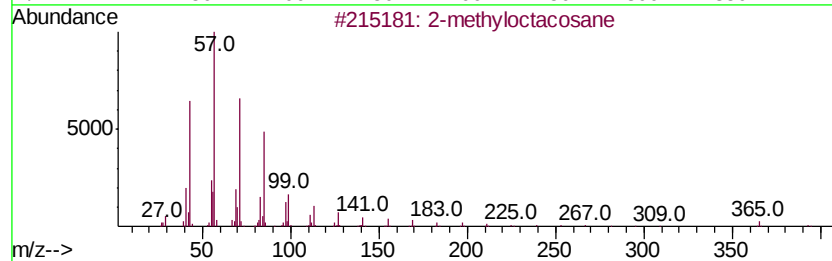
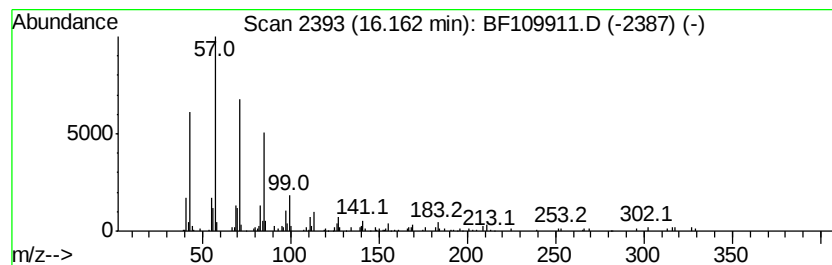
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 2-methyloctacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	2.59 ng	229029	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	86
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	86
3		Octacosane	394	C28H58	000630-02-4	86
4		2-Bromotetradecane	276	C14H29Br	074036-95-6	86
5		Tridecane, 3-methyl-	198	C14H30	006418-41-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\
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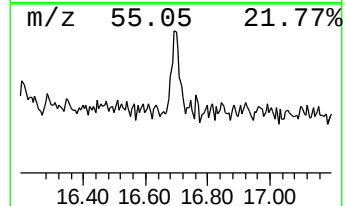
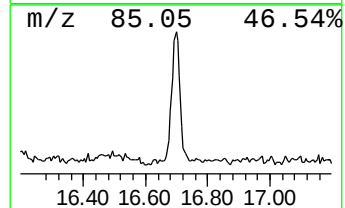
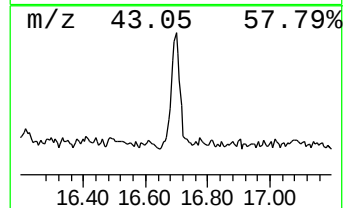
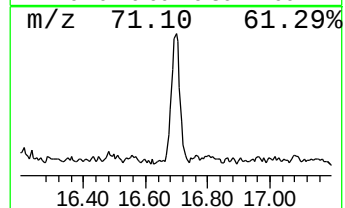
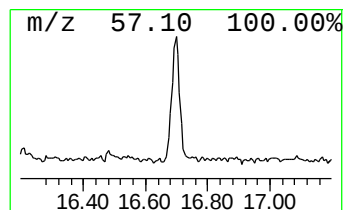
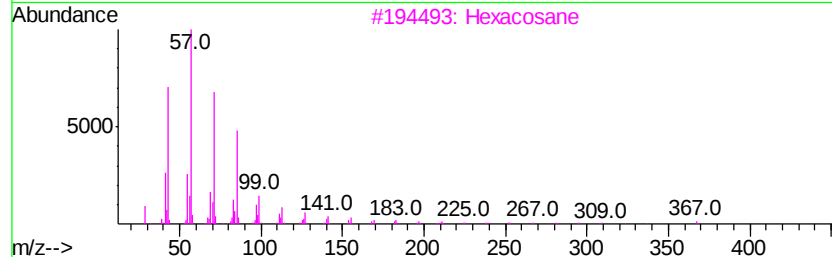
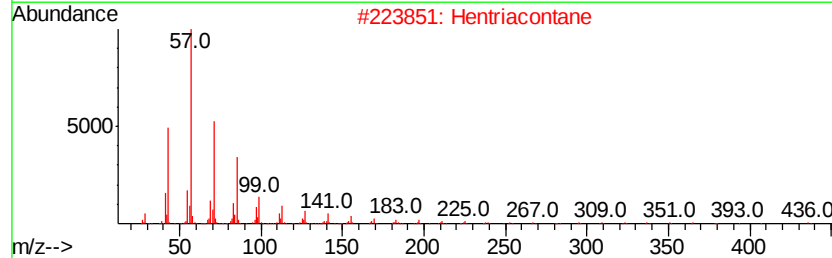
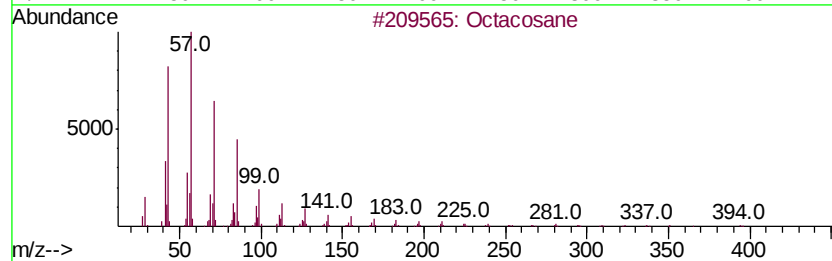
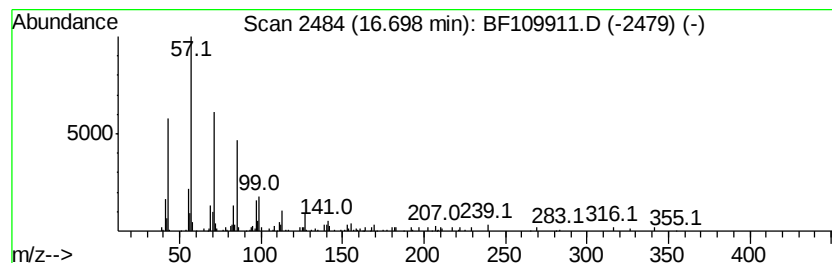
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Peak Number 13 Hentriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	2.17 ng	192229	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Hentriacontane	437	C31H64	000630-04-6	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Heneicosane	296	C21H44	000629-94-7	90



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 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1012-S-DH-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.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
Cyclohexane	2.12	6.8	ng	416261	1	6.98	1227630	20.0
Butane, 2-methoxy...	2.32	12.2	ng	746390	1	6.98	1227630	20.0
2-Pentanone, 4-hy...	5.21	2.2	ng	135874	1	6.98	1227630	20.0
unknown6.75	6.75	51.9	ng	3182980	1	6.98	1227630	20.0
9-Octadecenamide,...	13.57	7.1	ng	528566	5	14.16	1496390	20.0
7-Nonenamide	13.60	2.1	ng	157459	5	14.16	1496390	20.0
Octacosanol	13.97	2.9	ng	213352	5	14.16	1496390	20.0
Hexadecane	14.60	2.0	ng	153679	5	14.16	1496390	20.0
Octacosane	14.93	2.5	ng	217655	6	15.69	1769650	20.0
Heneicosane	15.29	2.6	ng	234794	6	15.69	1769650	20.0
2-methyloctacosane	16.16	2.6	ng	229029	6	15.69	1769650	20.0
Hentriacontane	16.70	2.2	ng	192229	6	15.69	1769650	20.0