

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
 Data File : BF113232.D
 Acq On : 22 Mar 2019 14:03
 Operator : JU/SJ
 Sample : K2040-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S4B(2-10)

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-BF022719.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	2.269	21	31	40	rVB2	49315	185863	3.25%	0.393%
2	5.340	547	553	556	rBV	3492229	3816301	66.78%	8.079%
3	6.369	722	728	744	rVB	3678118	4418106	77.31%	9.353%
4	6.492	744	749	752	rBV	4130028	4117801	72.05%	8.717%
5	6.716	783	787	790	rBV	1145434	985365	17.24%	2.086%
6	6.875	809	814	817	rBV	3336081	3259073	57.03%	6.899%
7	7.281	877	883	894	rBV	2349261	2757099	48.24%	5.837%
8	7.998	1000	1005	1011	rBV	1283408	1238104	21.66%	2.621%
9	9.075	1183	1188	1191	rBV	5241980	4971624	86.99%	10.525%
10	9.198	1206	1209	1213	rVB	200262	185317	3.24%	0.392%
11	9.463	1251	1254	1270	rVB	453138	438368	7.67%	0.928%
12	9.745	1298	1302	1313	rVB	1532446	1509567	26.41%	3.196%
13	10.339	1400	1403	1409	rBV	48393	58178	1.02%	0.123%
14	10.539	1432	1437	1440	rBV	3489709	3483571	60.95%	7.375%
15	11.227	1550	1554	1561	rBV	1917861	1706248	29.86%	3.612%
16	11.474	1592	1596	1603	rBV4	44483	76502	1.34%	0.162%
17	11.769	1641	1646	1654	rBV	888859	939884	16.45%	1.990%
18	12.280	1730	1733	1739	rBV	630842	943234	16.50%	1.997%
19	12.439	1754	1760	1761	rBV2	54657	62816	1.10%	0.133%
20	12.474	1762	1766	1774	rVB	202014	268890	4.70%	0.569%
21	12.551	1775	1779	1790	rBV	302647	431563	7.55%	0.914%
22	12.816	1819	1824	1827	rBV	5183379	5715053	100.00%	12.098%
23	13.051	1858	1864	1867	rBV2	162608	261171	4.57%	0.553%
24	13.498	1937	1940	1949	rVB4	48801	88819	1.55%	0.188%
25	13.721	1971	1978	1992	rBV2	413347	763395	13.36%	1.616%
26	13.857	1996	2001	2005	rBV	1600779	1728126	30.24%	3.658%
27	14.033	2028	2031	2038	rVB	88178	93206	1.63%	0.197%
28	14.339	2080	2083	2085	rBV	108208	100486	1.76%	0.213%
29	14.492	2105	2109	2112	rBV	100751	113201	1.98%	0.240%
30	14.633	2130	2133	2137	rVB	163038	151456	2.65%	0.321%
31	14.698	2141	2144	2148	rBV2	143609	149973	2.62%	0.317%
32	14.945	2183	2186	2193	rVB	223501	245012	4.29%	0.519%
33	15.268	2236	2241	2245	rBV	961947	1273796	22.29%	2.697%
34	15.298	2245	2246	2255	rVB	222044	225572	3.95%	0.478%

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Instrument :
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ClientSampleId :
S4B(2-10)

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.704	2310	2315	2320	rBV	199141	278683	4.88%	0.590%
36	16.168	2390	2394	2406	rVB4	106127	196338	3.44%	0.416%

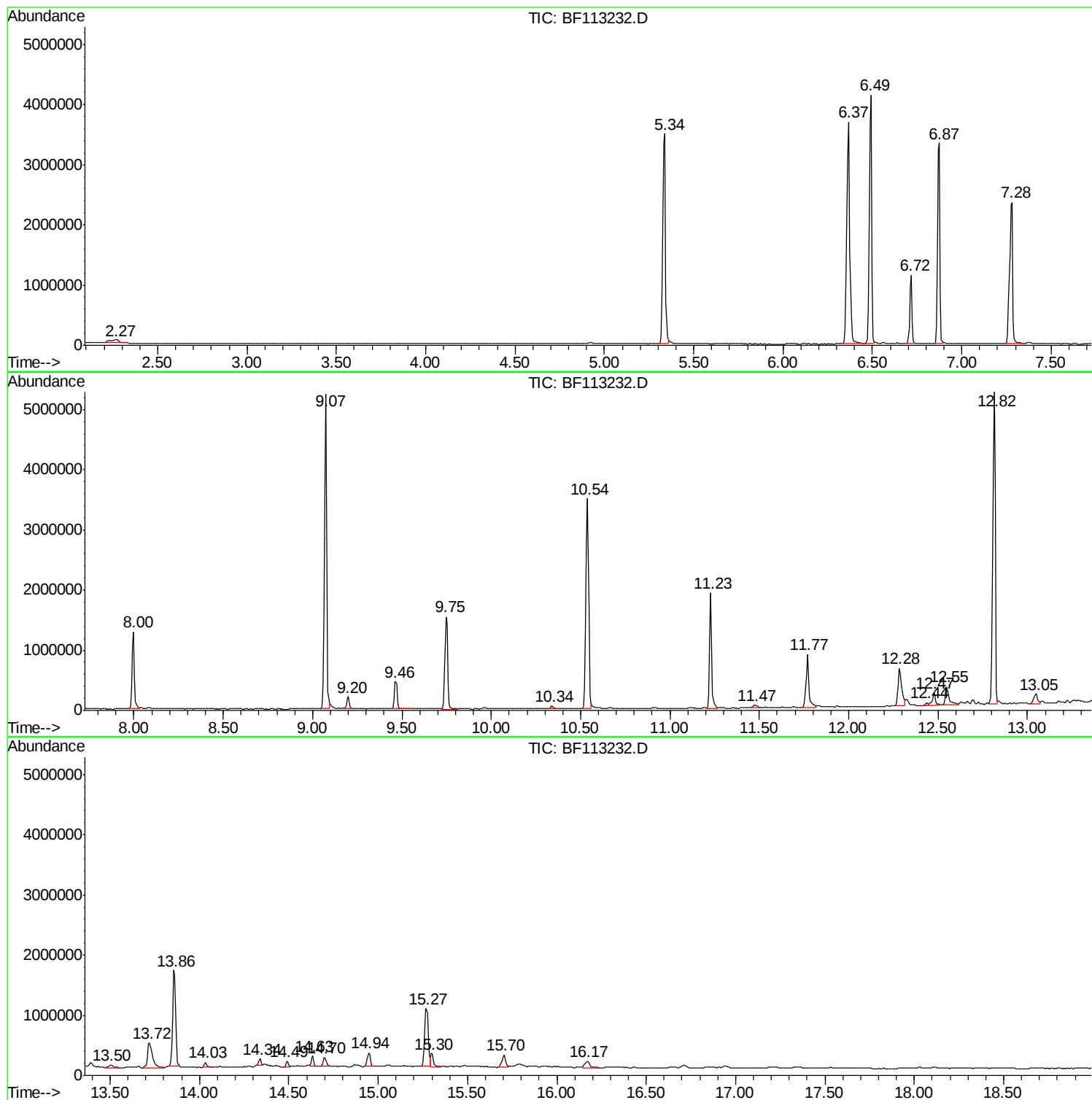
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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S4B(2-10)

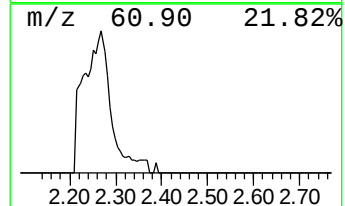
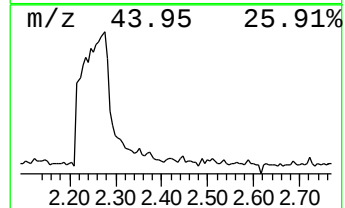
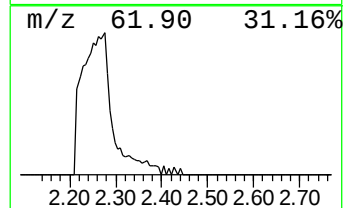
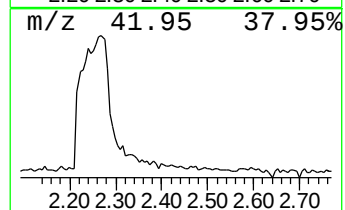
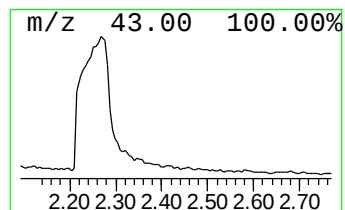
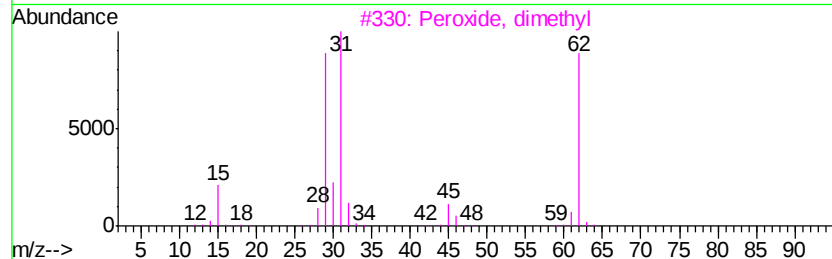
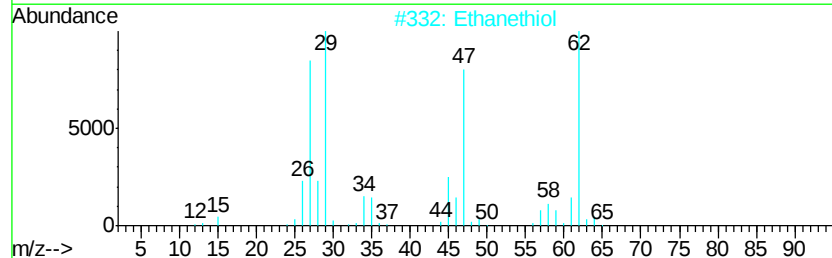
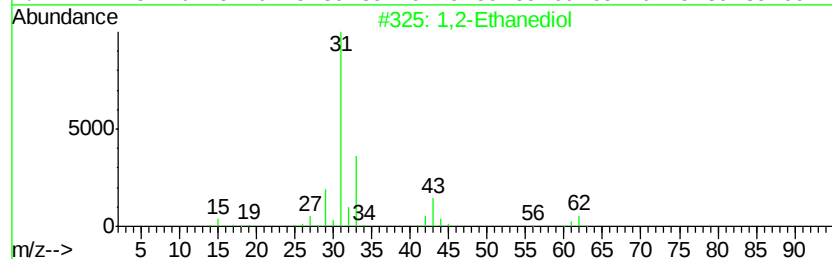
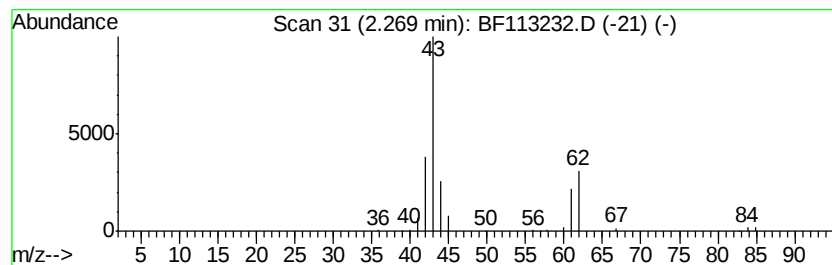
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.27 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	3.77 ng	185863	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Ethanediol	62	C2H6O2	000107-21-1	7
2		Ethanethiol	62	C2H6S	000075-08-1	4
3		Peroxide, dimethyl	62	C2H6O2	000690-02-8	4
4		Ethene, chloro-	62	C2H3Cl	000075-01-4	4
5		Pent-4-enylamine	85	C5H11N	022537-07-1	4



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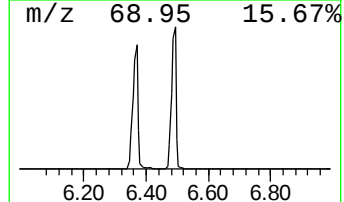
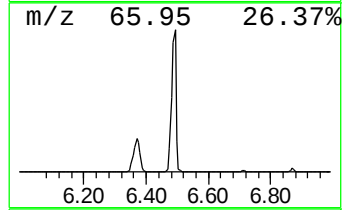
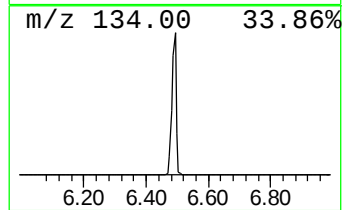
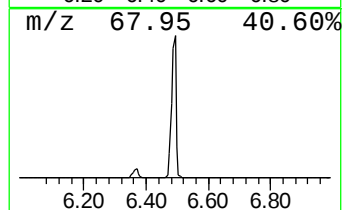
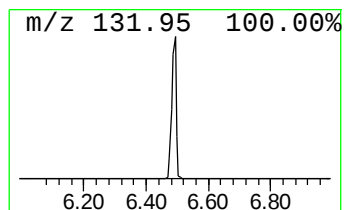
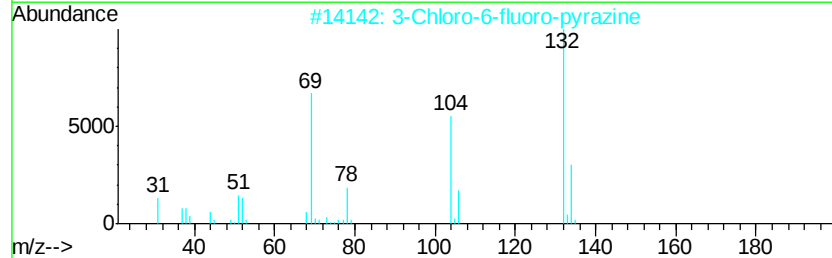
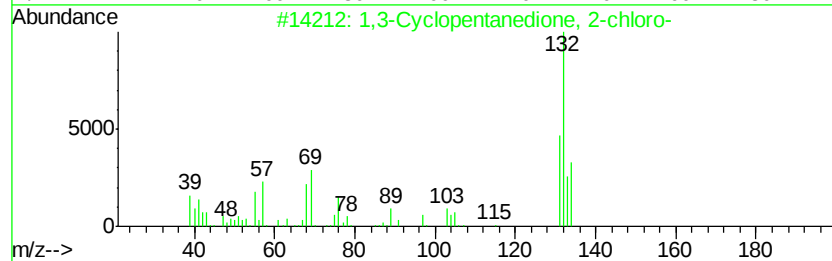
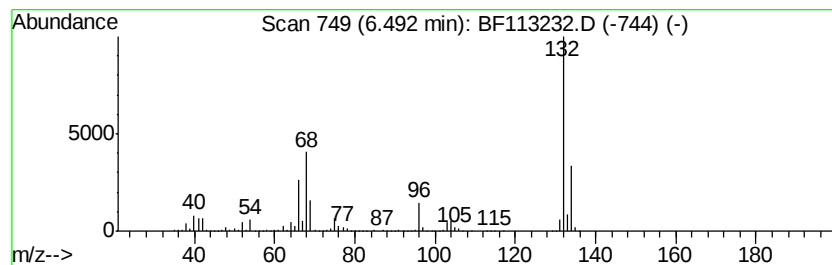
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	83.58 ng	4117800	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	47
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Xenon	132	Xe	007440-63-3	9



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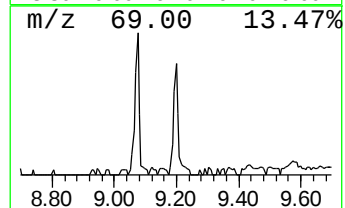
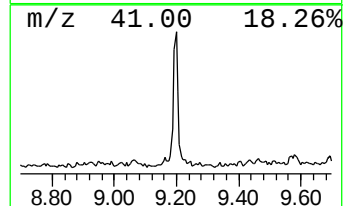
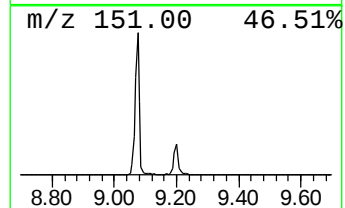
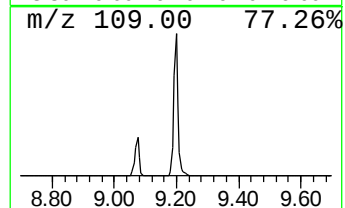
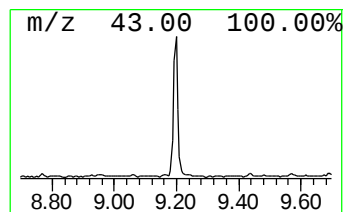
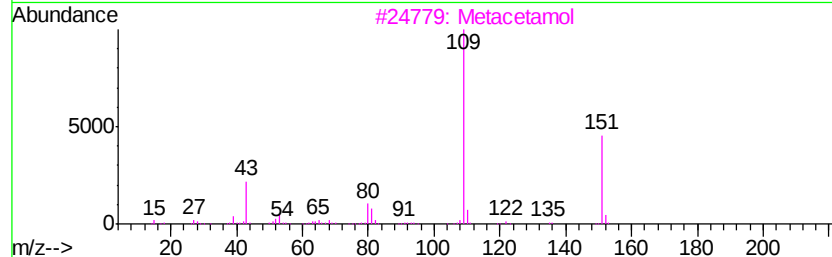
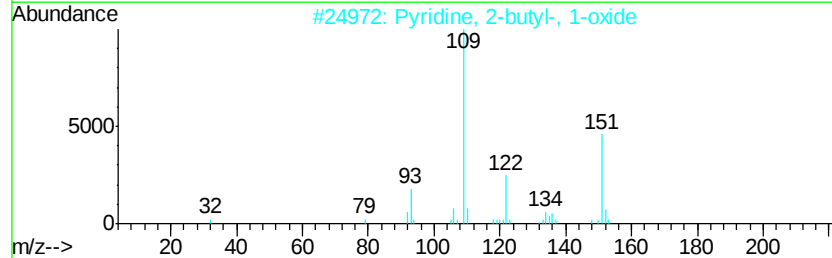
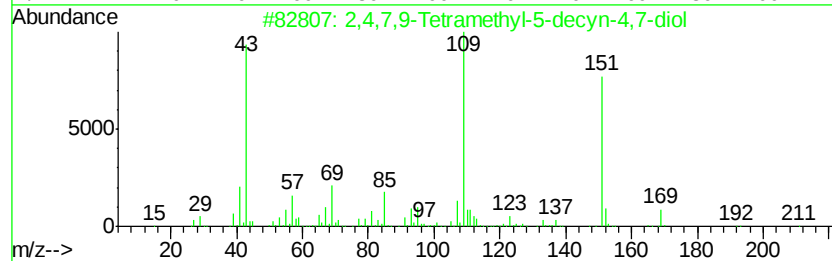
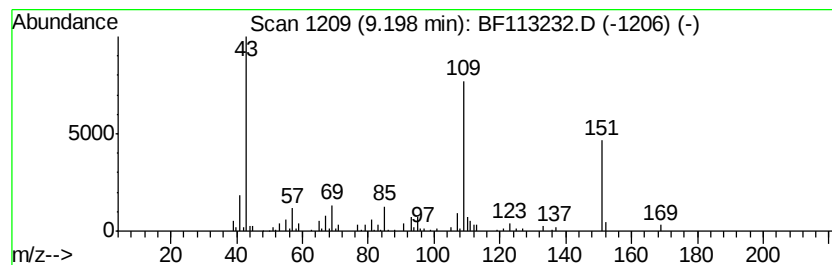
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.20	2.46 ng	185317	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2		Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
3		Metacetamol	151	C8H9NO2	000621-42-1	59
4		m-Isopropoxyaniline	151	C9H13NO	041406-00-2	59
5		Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	47



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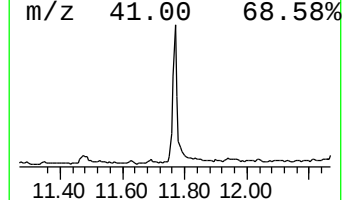
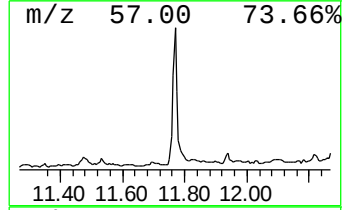
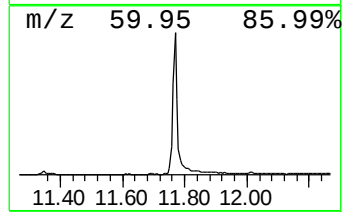
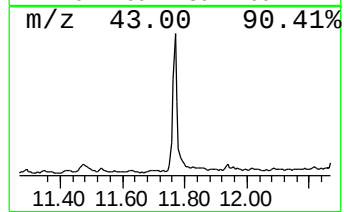
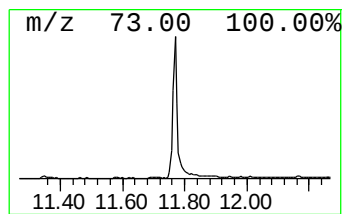
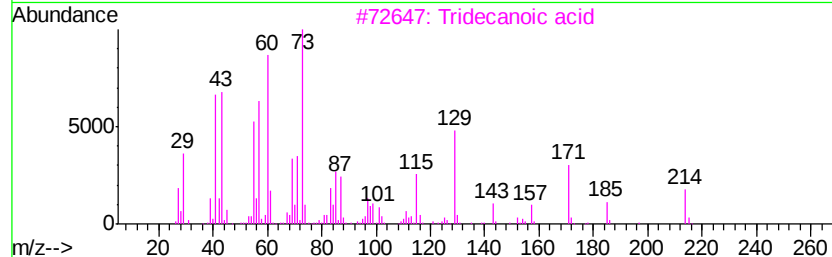
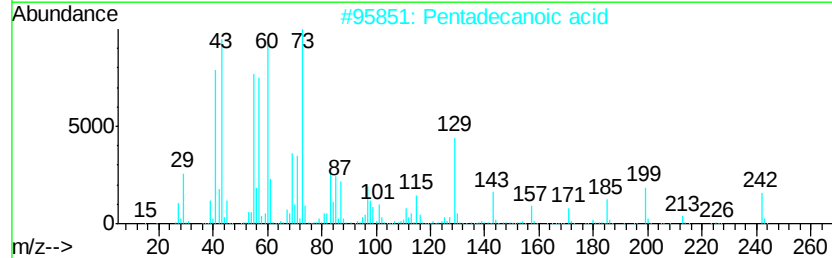
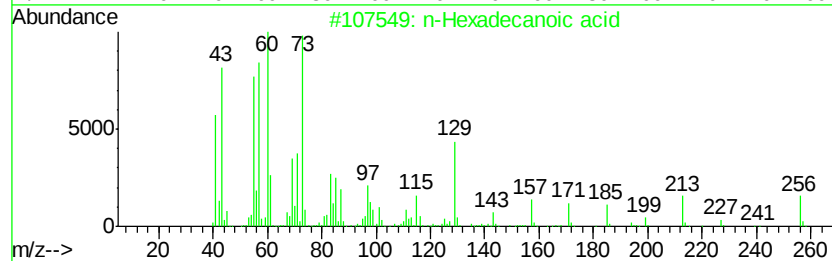
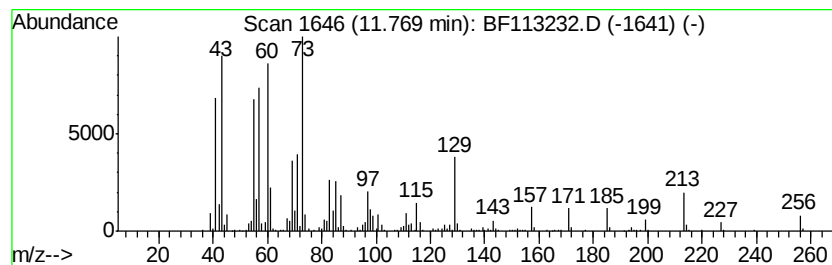
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	11.02 ng	939884	Phenanthrene-d10	11.23

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	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	92
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	81
5		n-Decanoic acid	172	C10H20O2	000334-48-5	62



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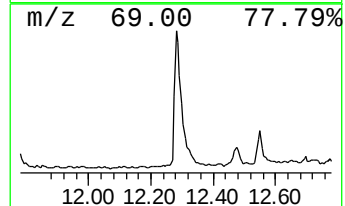
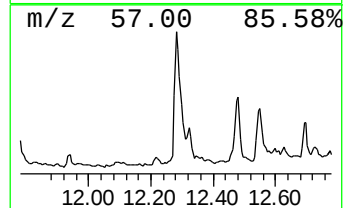
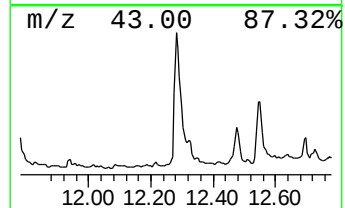
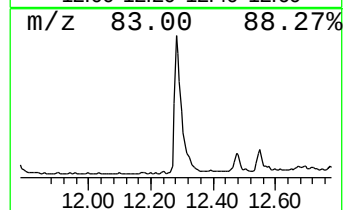
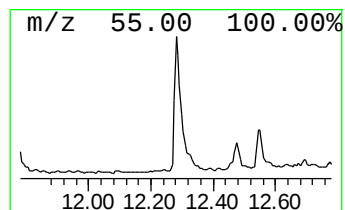
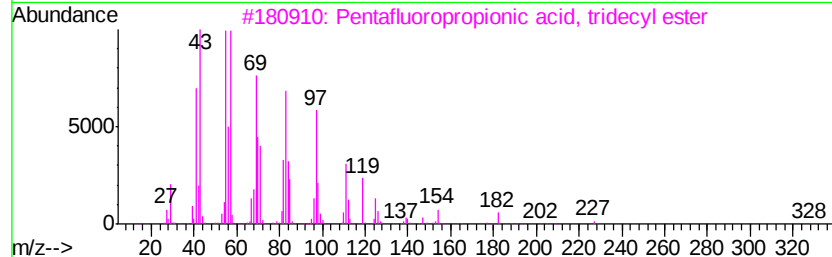
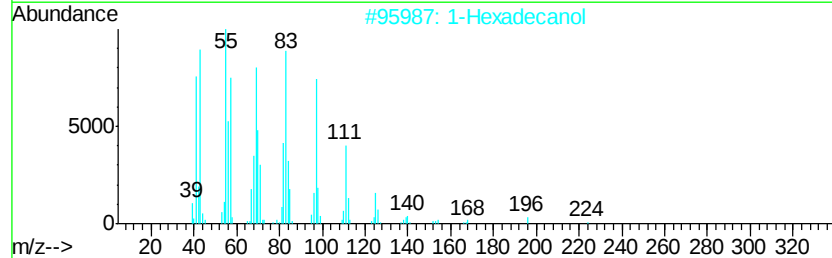
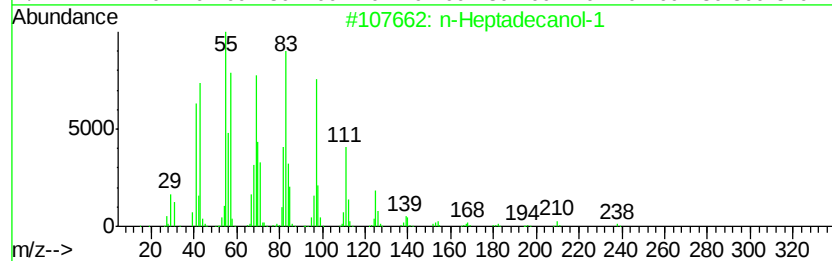
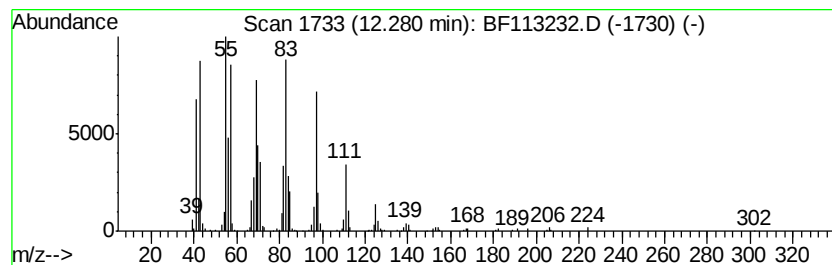
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Heptadecanol-1 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	11.06 ng	943234	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Heptadecanol-1	256	C17H36O	001454-85-9	95
2		1-Hexadecanol	242	C16H34O	036653-82-4	94
3		Pentafluoropropionic acid, tridecyl ester	346	C16H27F5O2	959261-22-4	93
4		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	91
5		Cyclohexadecane	224	C16H32	000295-65-8	91



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ALS Vial : 5 Sample Multiplier: 1

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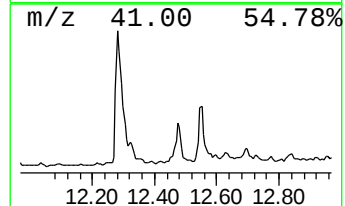
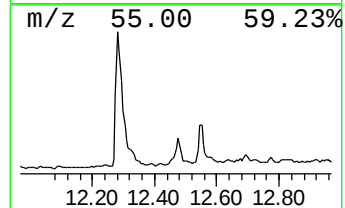
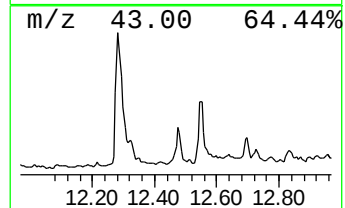
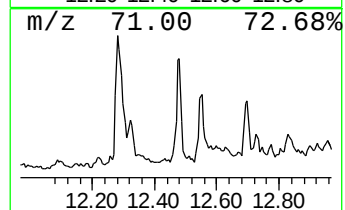
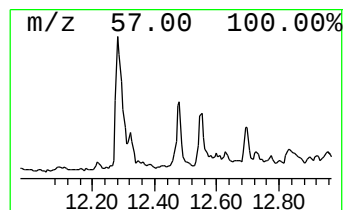
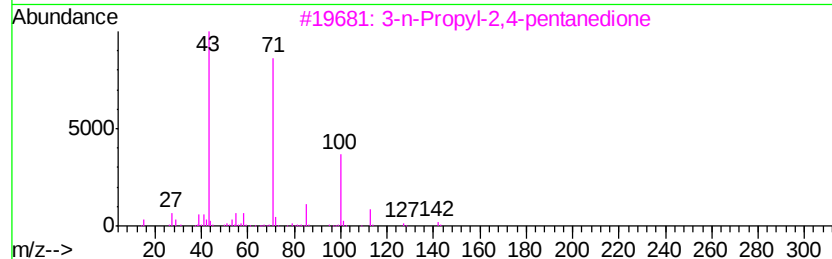
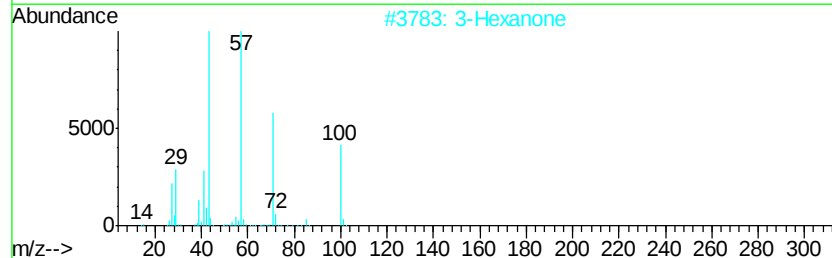
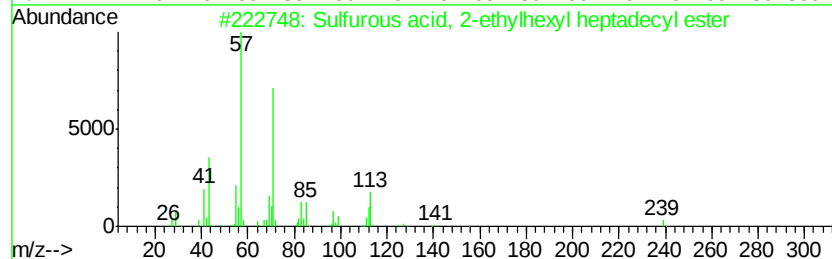
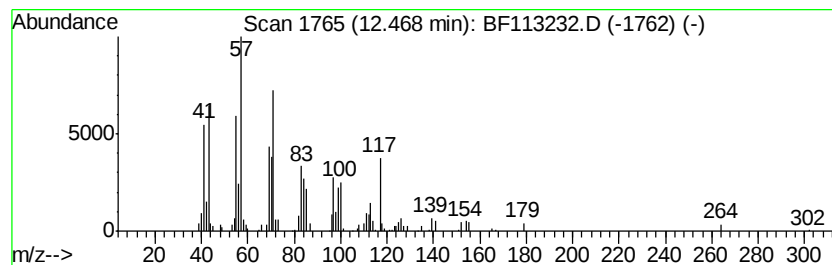
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown12.47 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	3.15 ng	268890	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	43
2		3-Hexanone	100	C6H12O	000589-38-8	38
3		3-n-Propyl-2,4-pentanedione	142	C8H14O2	001540-35-8	38
4		2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	38
5		Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
Operator : JU/SJ
Sample : K2040-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4B(2-10)

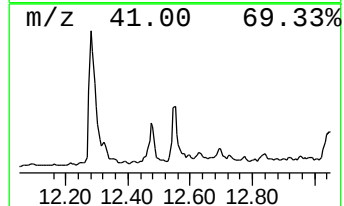
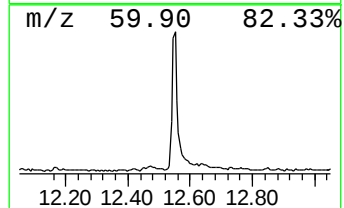
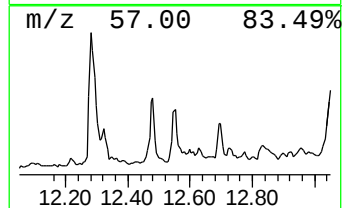
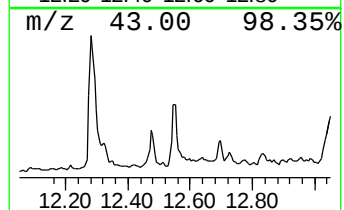
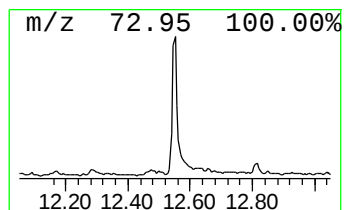
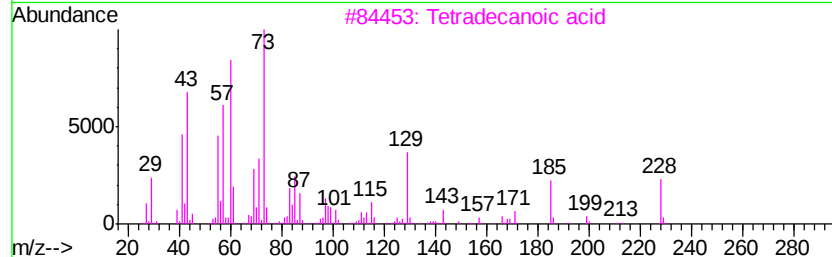
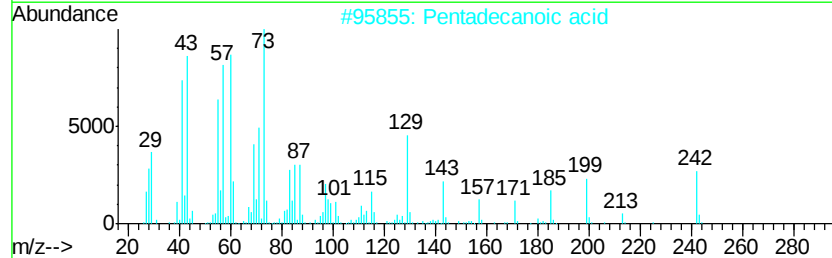
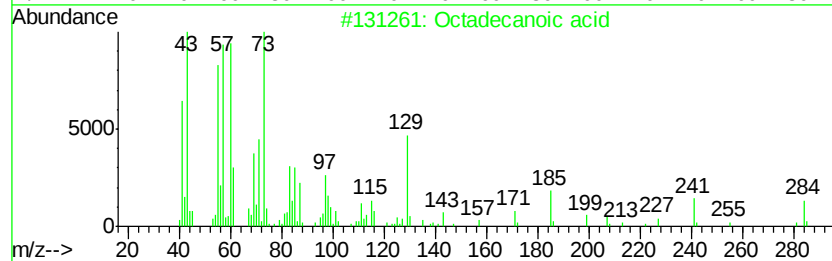
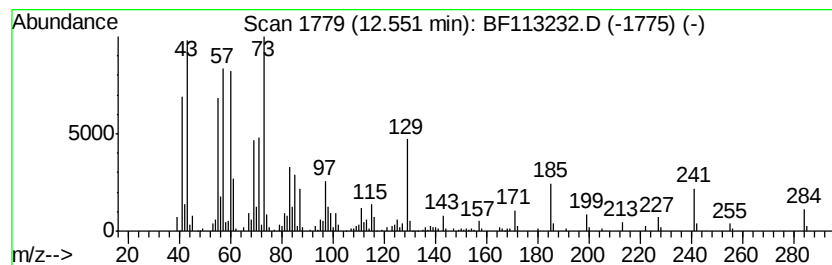
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	4.99 ng	431563	Chrysene-d12	13.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4		Undecanoic acid	186	C11H22O2	000112-37-8	72
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
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Sample : K2040-01
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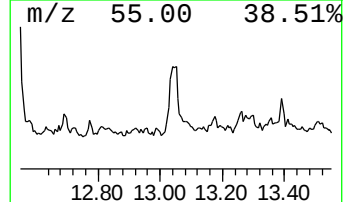
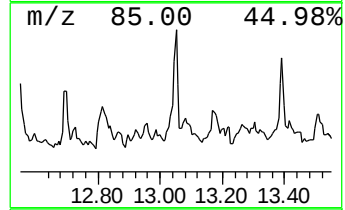
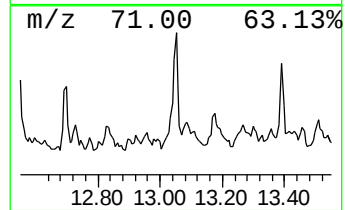
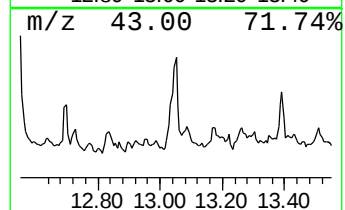
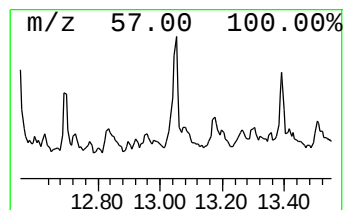
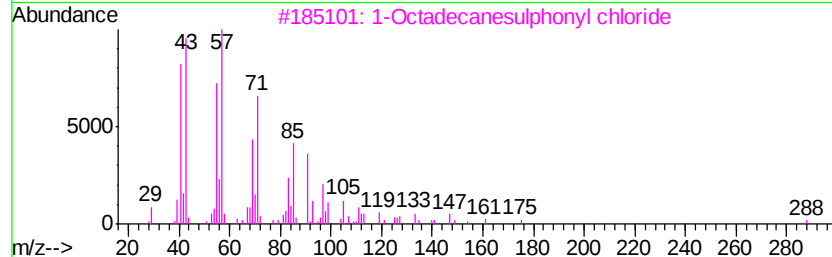
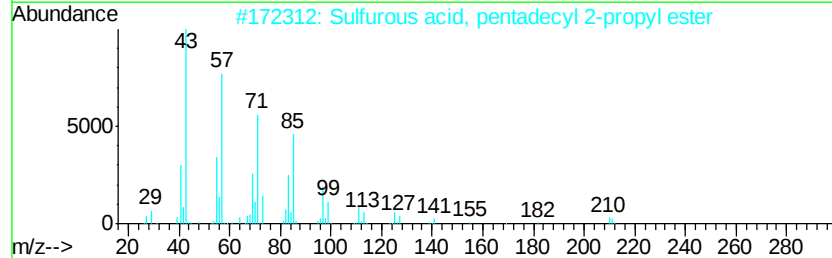
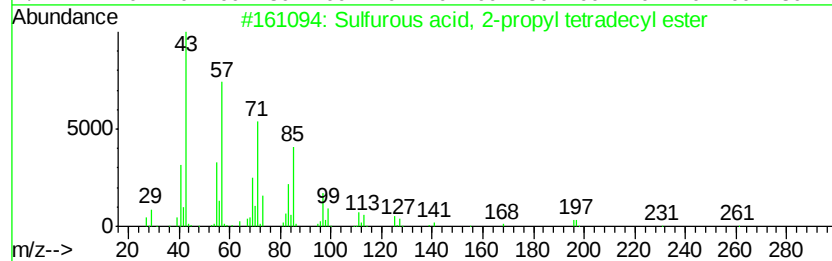
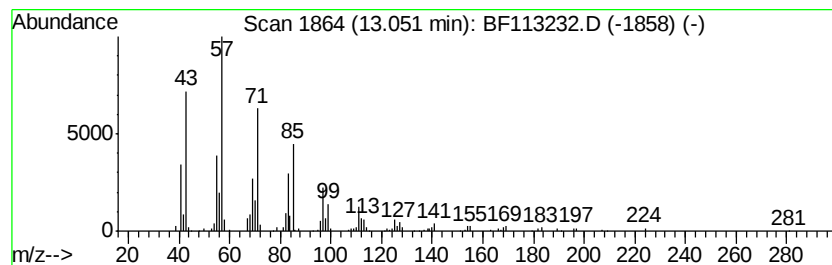
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Sulfurous acid, 2-propyl te... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.05	3.02 ng	261171	Chrysene-d12	13.86		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	91
2		Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	91
3		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	87
4		Tritetracontane	605	C43H88	007098-21-7	83
5		Tetratetracontane	619	C44H90	007098-22-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
Operator : JU/SJ
Sample : K2040-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4B(2-10)

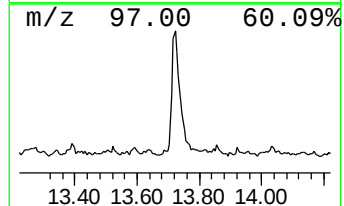
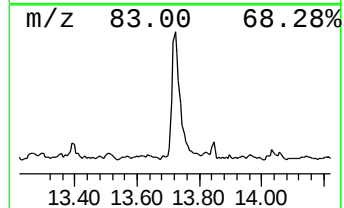
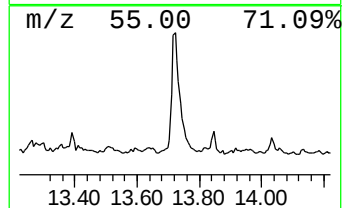
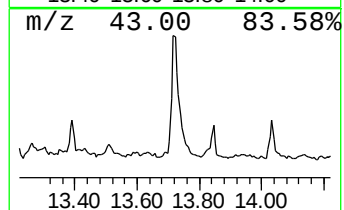
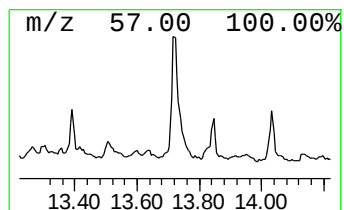
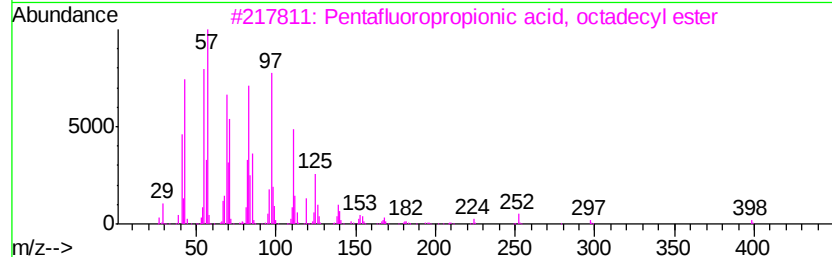
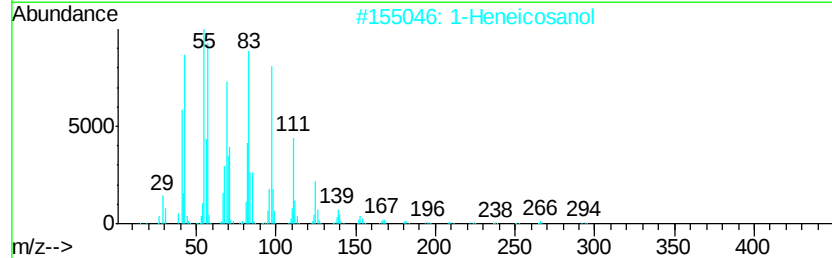
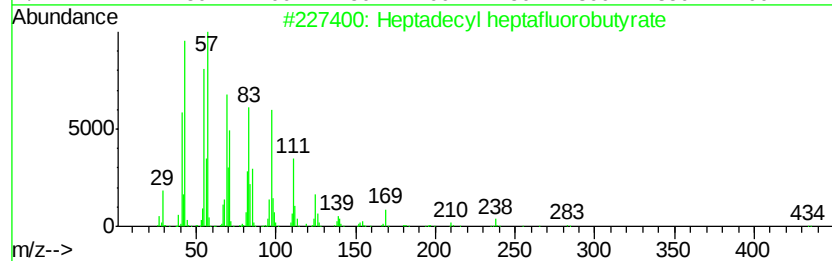
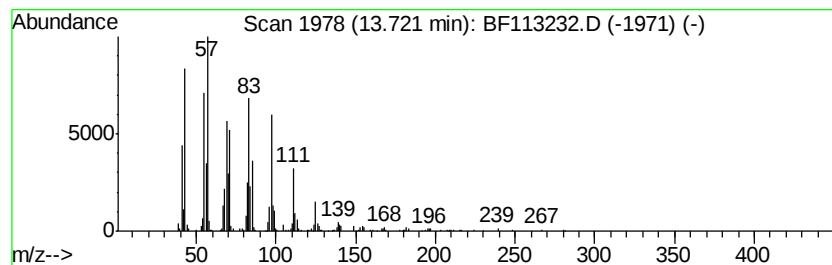
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecyl heptafluorobutyrate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	8.83 ng	763395	Chrysene-d12	13.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	91
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
5		Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
 Data File : BF113232.D
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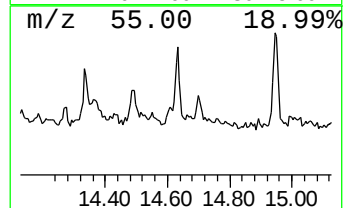
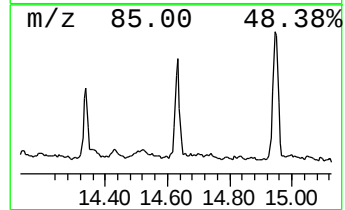
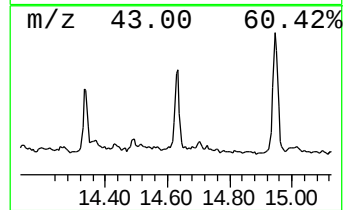
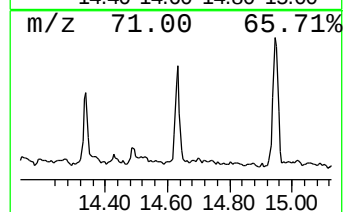
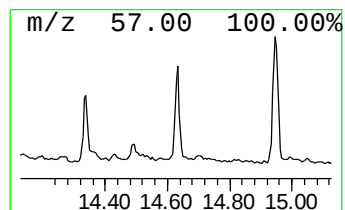
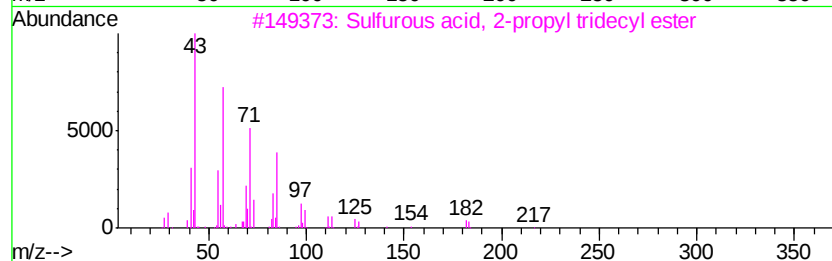
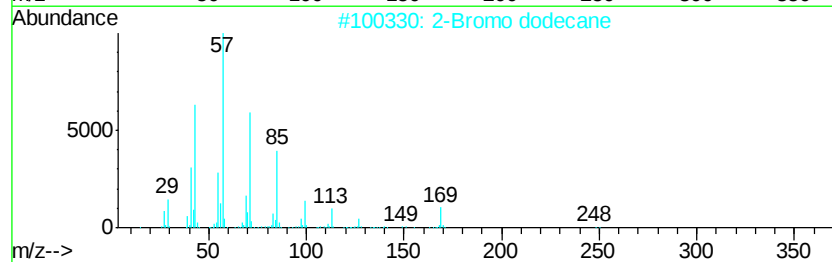
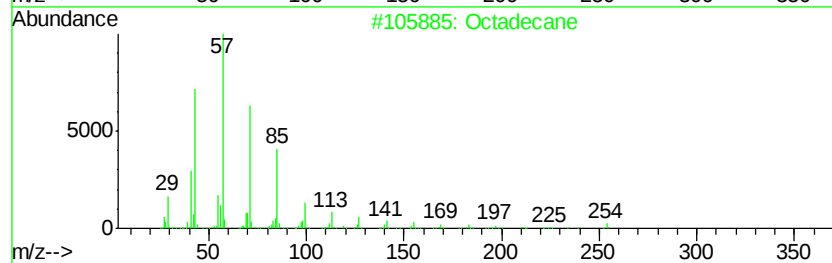
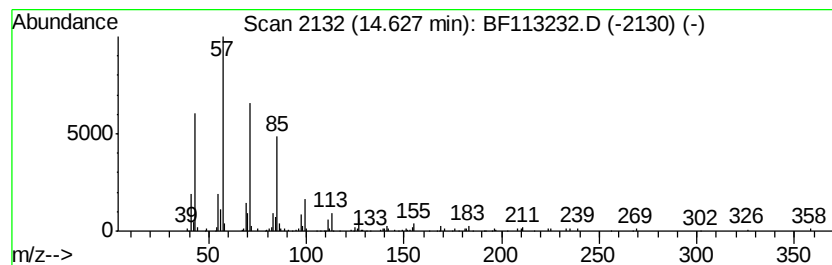
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Octadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	2.38 ng	151456	Perylene-d12	15.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	86
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	86
3		Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	86
4		Octacosane	394	C28H58	000630-02-4	86
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
Operator : JU/SJ
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S4B(2-10)

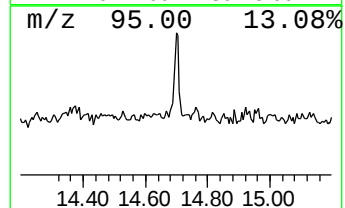
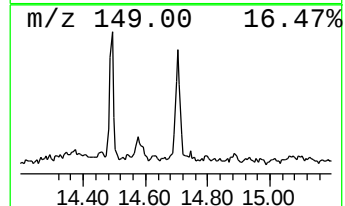
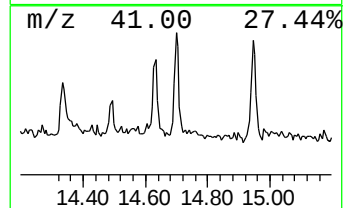
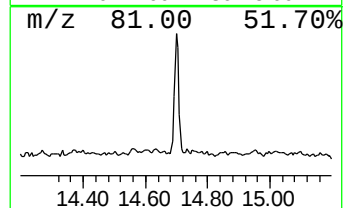
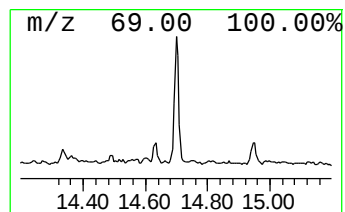
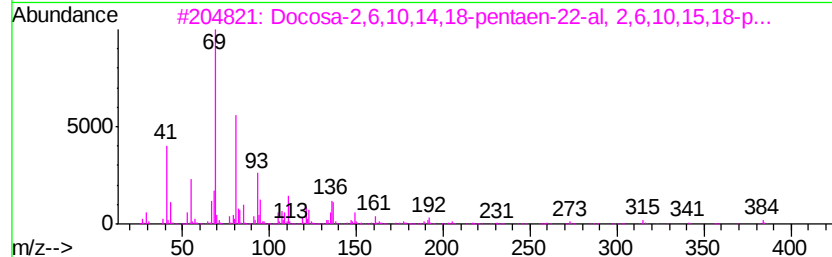
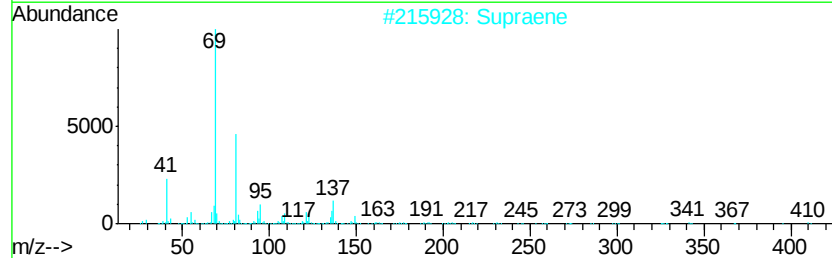
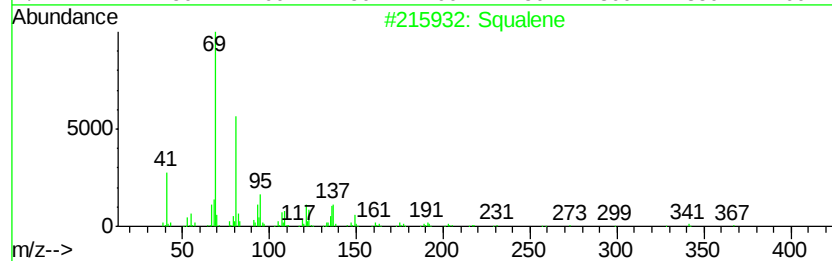
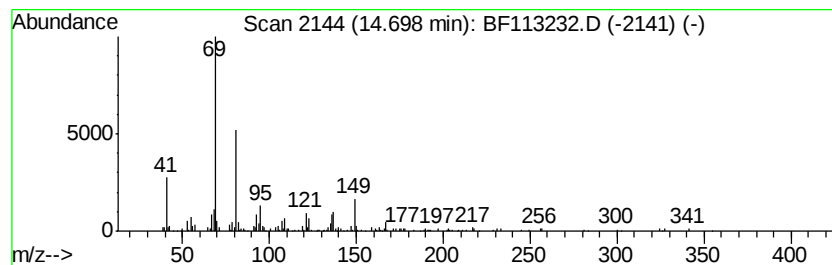
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	2.35 ng	149973	Perylene-d12	15.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	87
2		Supraene	410	C30H50	007683-64-9	74
3		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	72
4		Trideca-1,3,7,11-tetraene-1,1-di...	268	C18H24N2	1000159-88-9	53
5		Crotonic acid, o-formylphenyl ester	190	C11H10O3	114636-69-0	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
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Operator : JU/SJ
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Misc :
ALS Vial : 5 Sample Multiplier: 1

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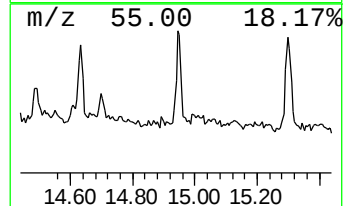
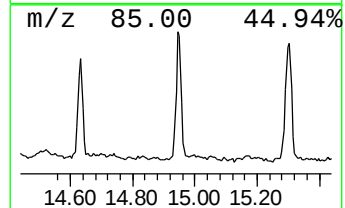
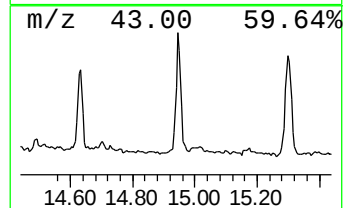
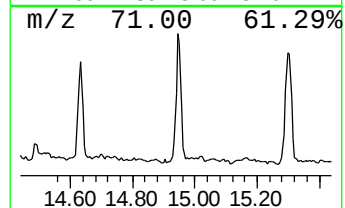
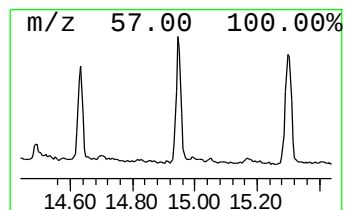
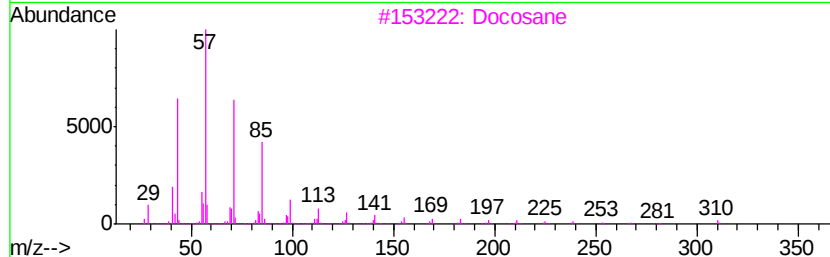
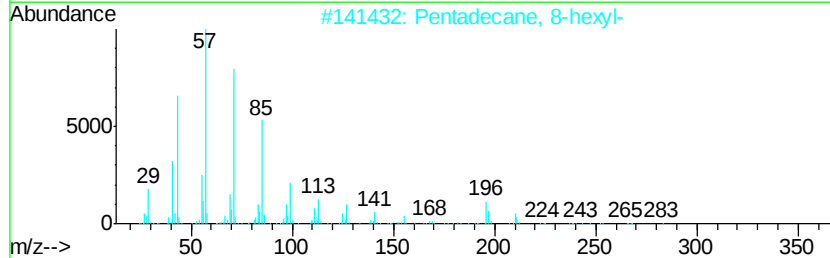
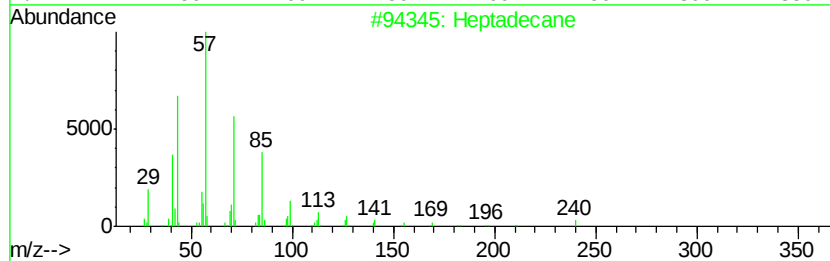
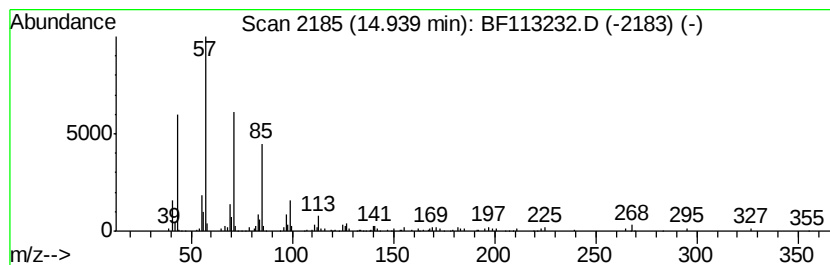
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	3.85 ng	245012	Perylene-d12	15.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	91
3	Docosane		310	C22H46	000629-97-0	91
4	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	90
5	Nonadecane		268	C19H40	000629-92-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
Operator : JU/SJ
Sample : K2040-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4B(2-10)

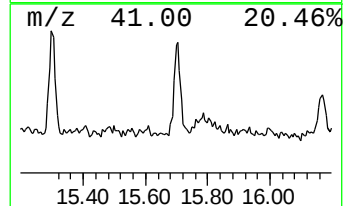
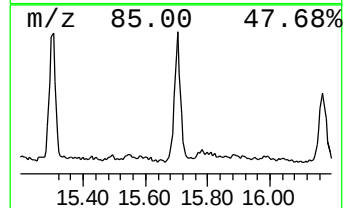
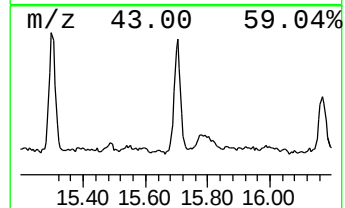
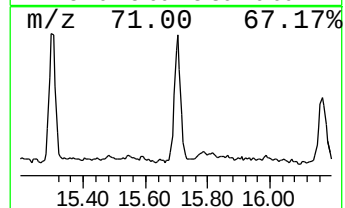
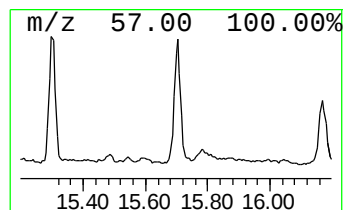
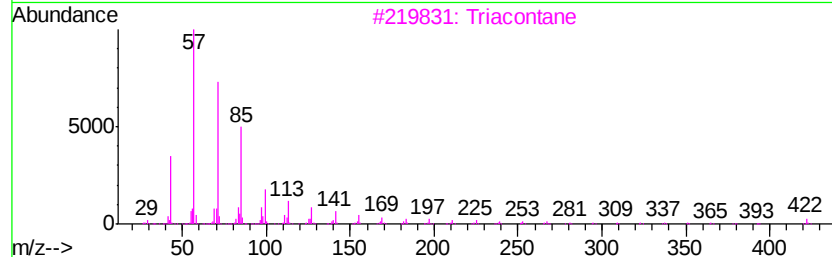
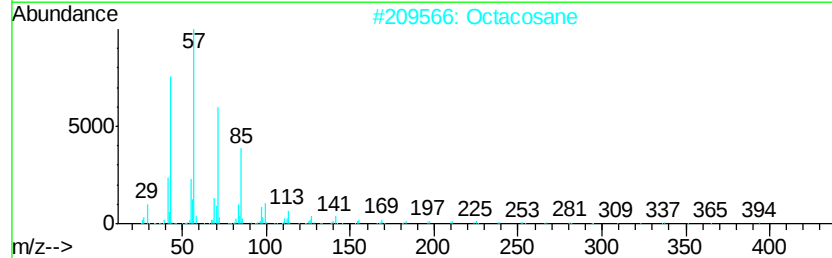
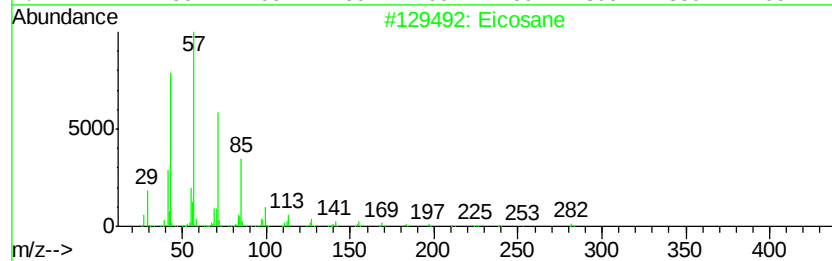
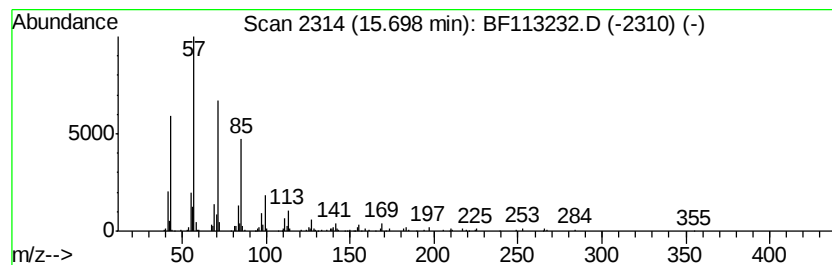
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	4.38 ng	278683	Perylene-d12	15.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Octacosane	394	C28H58	000630-02-4	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
Data File : BF113232.D
Acq On : 22 Mar 2019 14:03
Operator : JU/SJ
Sample : K2040-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4B(2-10)

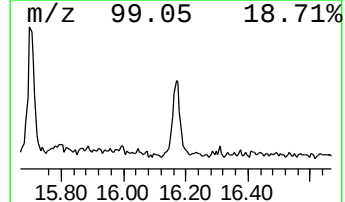
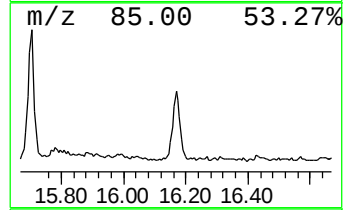
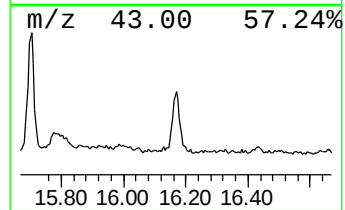
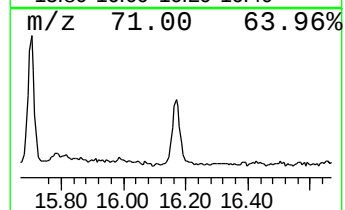
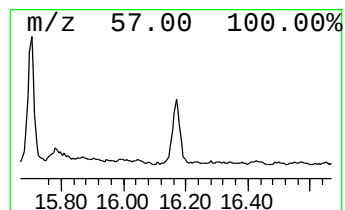
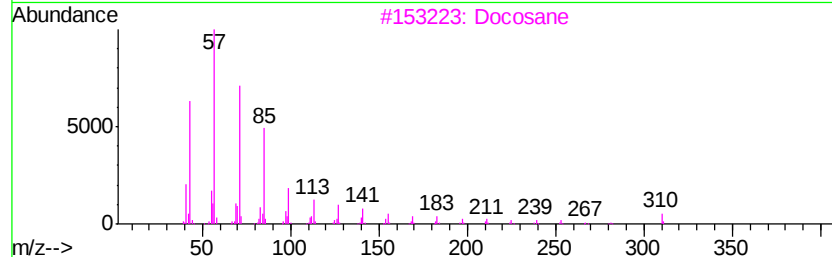
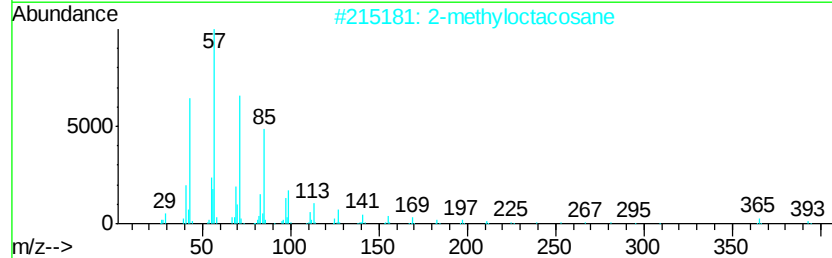
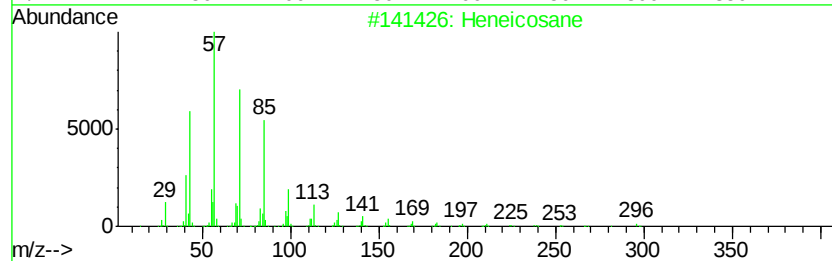
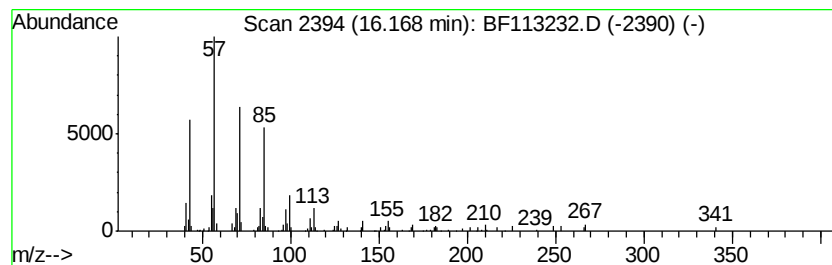
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.17	3.08 ng	196338	Perylene-d12	15.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Docosane	310	C22H46	000629-97-0	86
4		Tridecane, 3-methyl-	198	C14H30	006418-41-3	86
5		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
 Data File : BF113232.D
 Acq On : 22 Mar 2019 14:03
 Operator : JU/SJ
 Sample : K2040-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S4B(2-10)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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
unknown2.27	2.27	3.8	ng	185863	1	6.72	985365	20.0
unknown6.49	6.49	83.6	ng	4117800	1	6.72	985365	20.0
2,4,7,9-Tetrameth...	9.20	2.5	ng	185317	3	9.75	1509570	20.0
n-Hexadecanoic acid	11.77	11.0	ng	939884	4	11.23	1706250	20.0
n-Heptadecanol-1	12.28	11.1	ng	943234	4	11.23	1706250	20.0
unknown12.47	12.47	3.1	ng	268890	4	11.23	1706250	20.0
Octadecanoic acid	12.55	5.0	ng	431563	5	13.86	1728130	20.0
Sulfurous acid, 2...	13.05	3.0	ng	261171	5	13.86	1728130	20.0
Heptadecyl heptaf...	13.72	8.8	ng	763395	5	13.86	1728130	20.0
Octadecane	14.63	2.4	ng	151456	6	15.27	1273800	20.0
Squalene	14.70	2.4	ng	149973	6	15.27	1273800	20.0
Heptadecane	14.94	3.9	ng	245012	6	15.27	1273800	20.0
Eicosane	15.70	4.4	ng	278683	6	15.27	1273800	20.0
Heneicosane	16.17	3.1	ng	196338	6	15.27	1273800	20.0