

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113467.D
 Acq On : 1 Apr 2019 15:32
 Operator : JU/SJ
 Sample : K2171-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-COMP

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-BF032519.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	5.322	546	550	576	rBV	2410730	2580427	85.50%	8.366%
2	6.351	721	725	743	rBV	2395782	2723350	90.24%	8.830%
3	6.481	743	747	750	rBV	2927847	2682743	88.89%	8.698%
4	6.704	782	785	794	rBV	1019423	857524	28.41%	2.780%
5	6.863	808	812	829	rVB	2342639	2108446	69.86%	6.836%
6	7.269	877	881	904	rBV	1701260	1731249	57.36%	5.613%
7	7.986	999	1003	1018	rBV	1160163	1112795	36.87%	3.608%
8	9.063	1181	1186	1205	rBV	2930892	3018063	100.00%	9.785%
9	9.457	1249	1253	1262	rBV	219454	216184	7.16%	0.701%
10	9.733	1296	1300	1318	rVB	1285224	1268724	42.04%	4.113%
11	10.522	1430	1434	1448	rBV	1819801	1882531	62.38%	6.104%
12	11.216	1548	1552	1569	rBV	1376512	1370751	45.42%	4.444%
13	11.751	1639	1643	1653	rBV2	154215	231251	7.66%	0.750%
14	12.421	1754	1757	1761	rBV	43631	44933	1.49%	0.146%
15	12.527	1771	1775	1789	rBV	232465	403997	13.39%	1.310%
16	12.651	1793	1796	1799	rVB	46391	42230	1.40%	0.137%
17	12.792	1816	1820	1829	rBV	3060316	2846378	94.31%	9.229%
18	13.333	1908	1912	1915	rBV	74260	67657	2.24%	0.219%
19	13.698	1971	1974	1977	rBV	82981	77864	2.58%	0.252%
20	13.745	1977	1982	1993	rVB7	32105	96331	3.19%	0.312%
21	13.845	1993	1999	2007	rBV	1024995	1139074	37.74%	3.693%
22	14.010	2024	2027	2032	rBV	173138	166424	5.51%	0.540%
23	14.315	2075	2079	2083	rBV	316418	274317	9.09%	0.889%
24	14.474	2103	2106	2109	rBV4	23139	31208	1.03%	0.101%
25	14.610	2125	2129	2133	rVB	462207	427719	14.17%	1.387%
26	14.674	2137	2140	2146	rBV	196184	201654	6.68%	0.654%
27	14.857	2168	2171	2177	rBV2	23937	36636	1.21%	0.119%
28	14.921	2178	2182	2190	rVB	485104	516820	17.12%	1.676%
29	15.133	2215	2218	2221	rBV	37001	43710	1.45%	0.142%
30	15.251	2233	2238	2248	rBV2	869149	1646775	54.56%	5.339%
31	15.668	2303	2309	2316	rBV	351605	515611	17.08%	1.672%
32	16.127	2381	2387	2398	rVB	192661	353068	11.70%	1.145%
33	16.662	2474	2478	2485	rVB2	68475	126528	4.19%	0.410%

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ClientSampleId :
GW-COMP

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

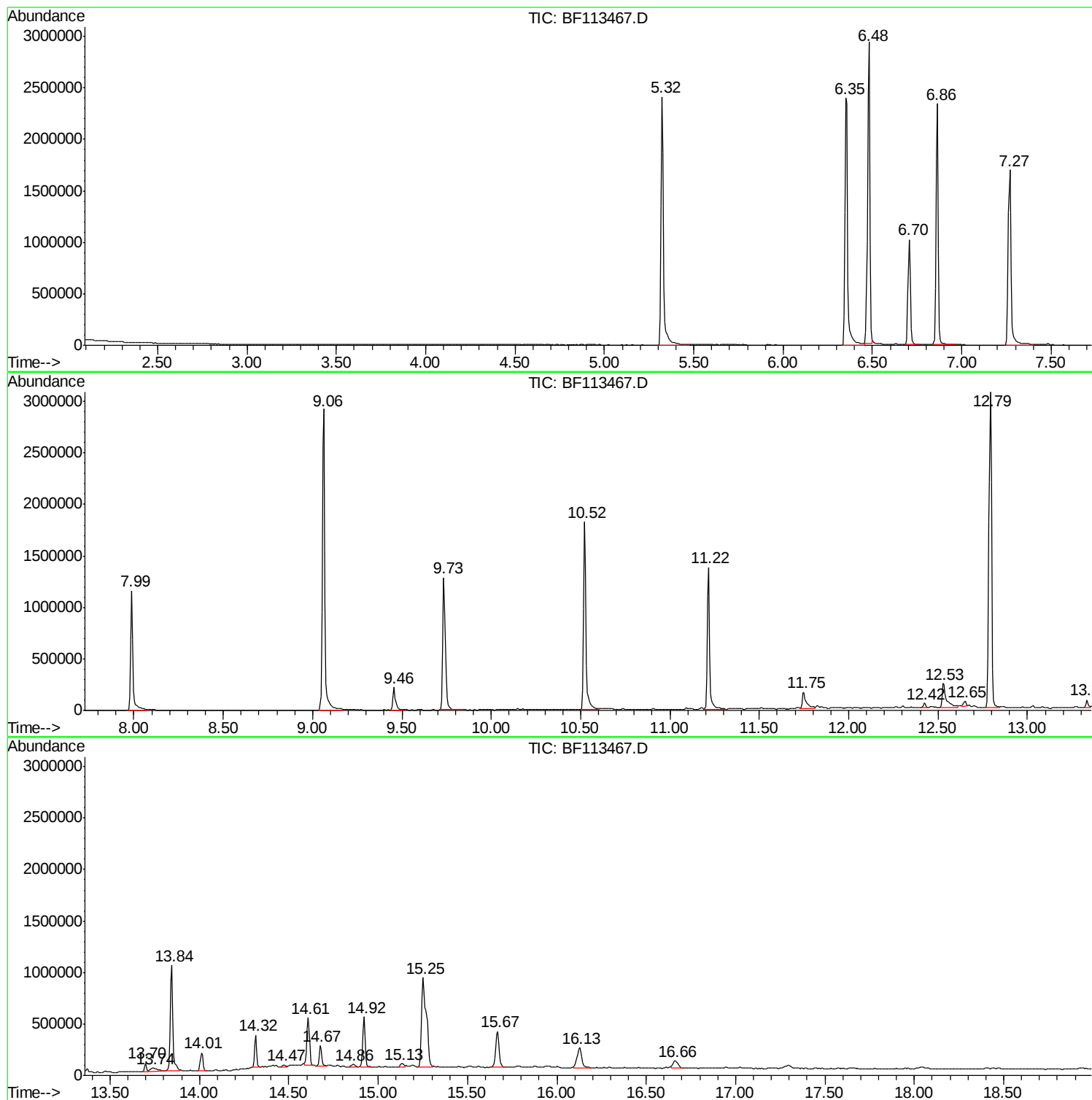
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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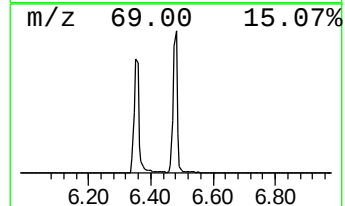
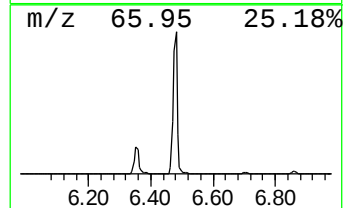
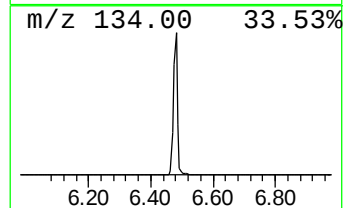
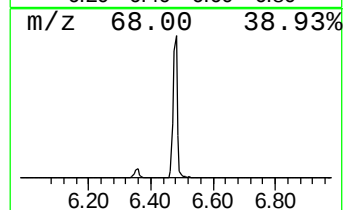
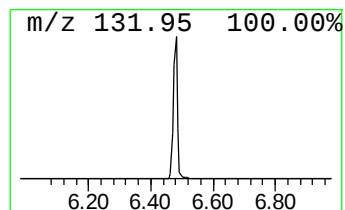
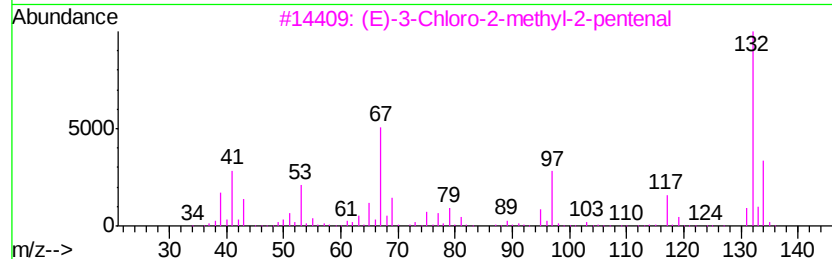
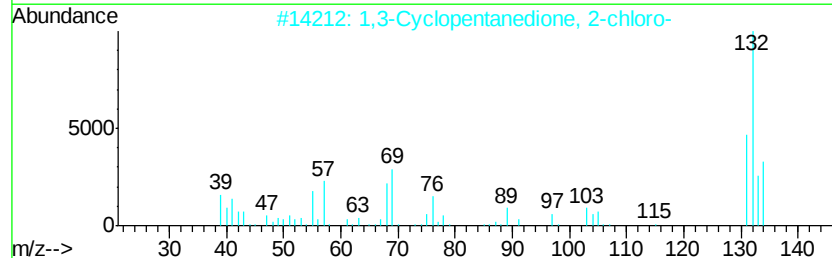
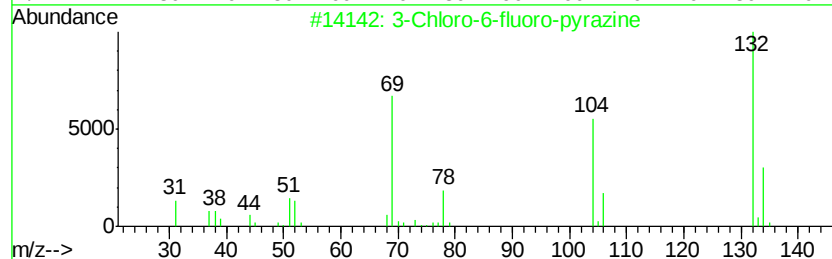
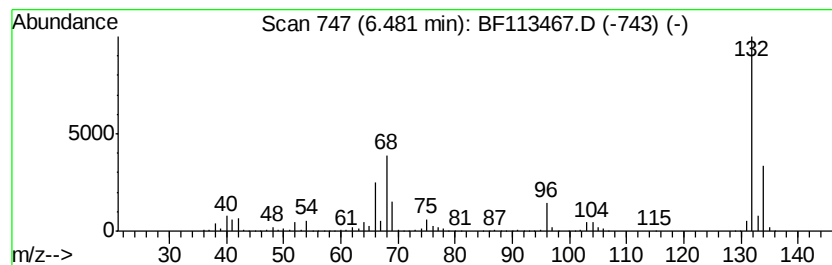
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	62.57 ng	2682740	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Xenon	132	Xe	007440-63-3	9



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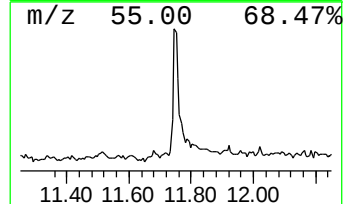
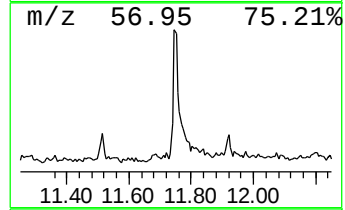
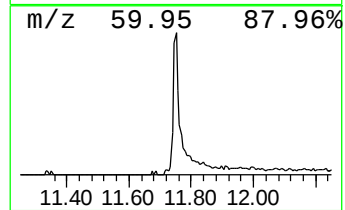
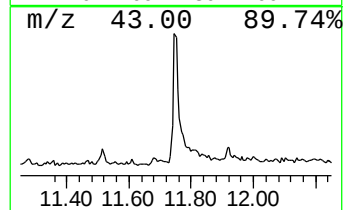
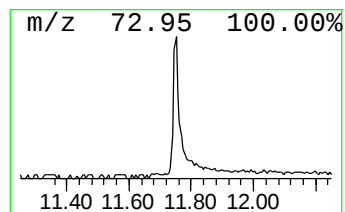
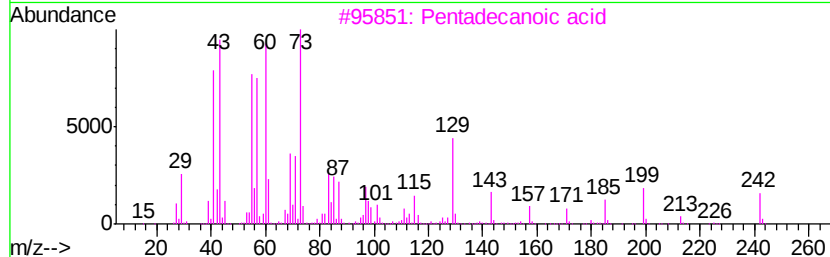
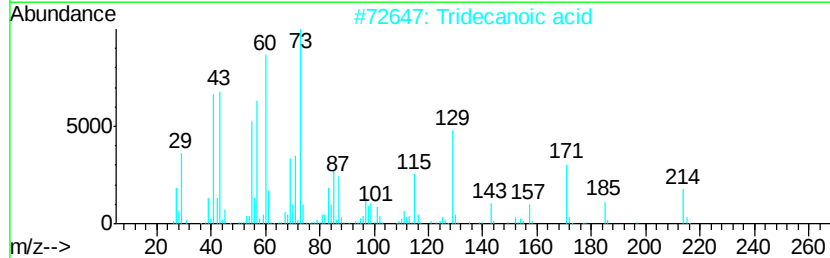
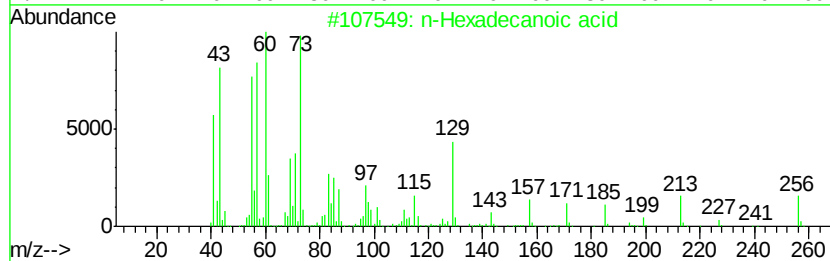
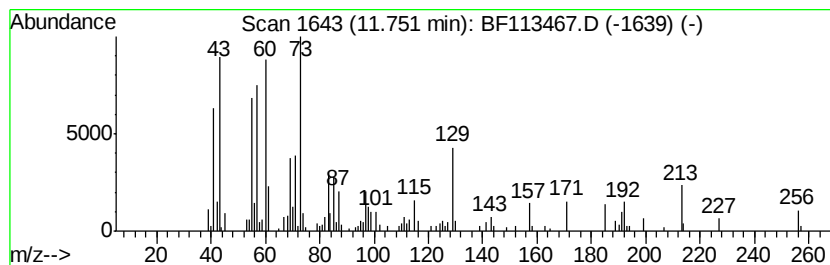
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.75	3.37 ng	231251	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	20



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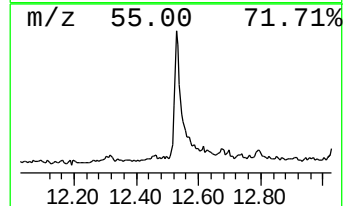
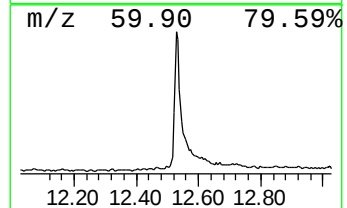
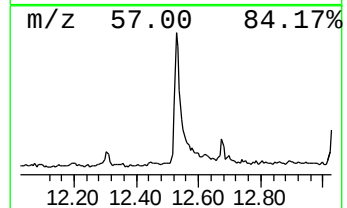
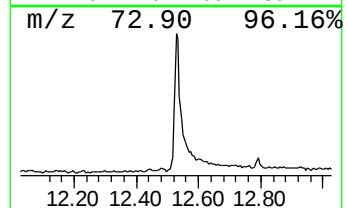
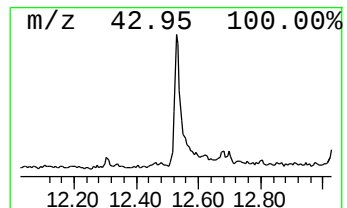
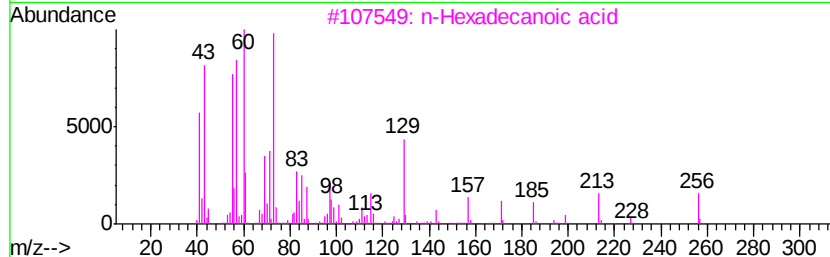
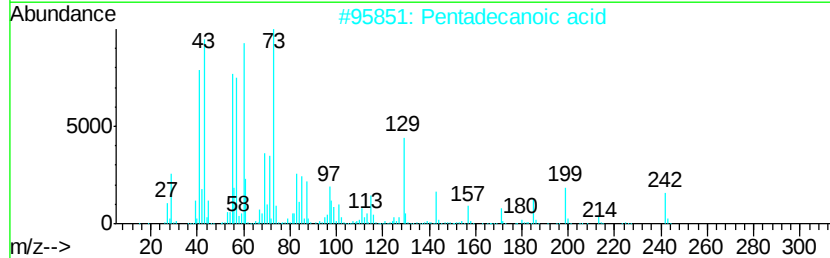
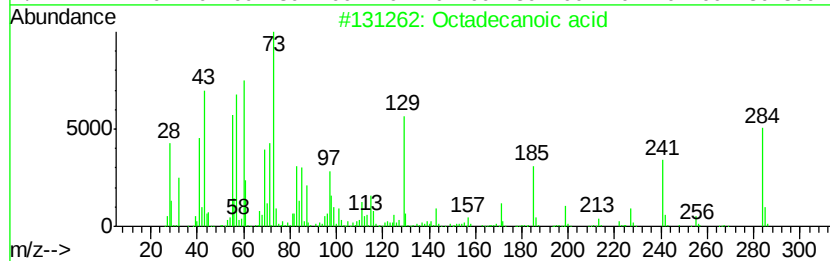
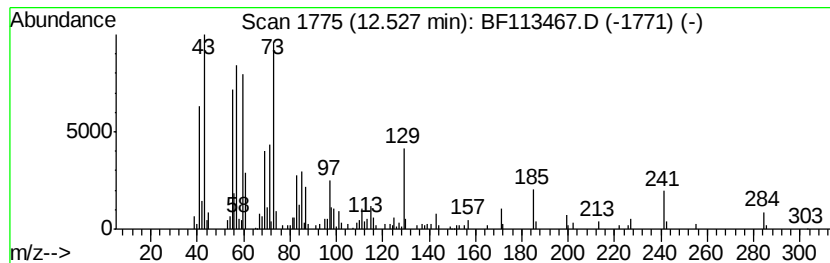
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	5.89 ng	403997	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	70



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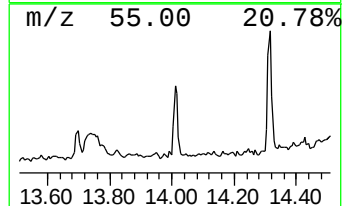
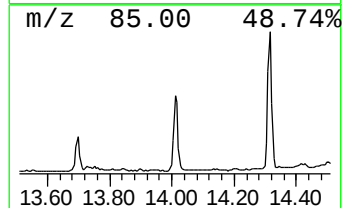
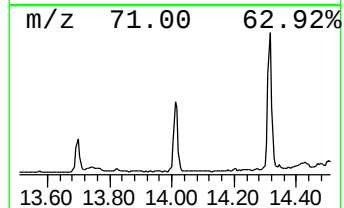
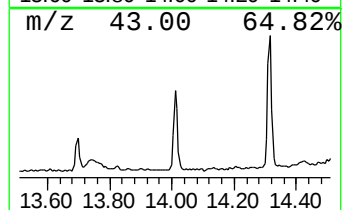
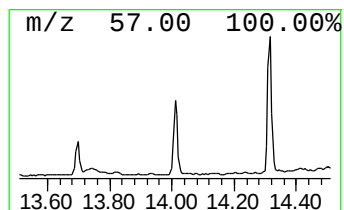
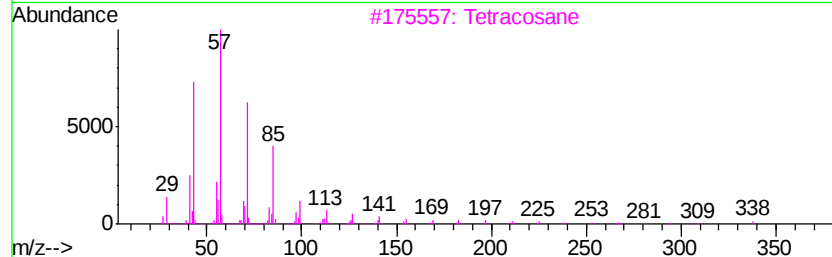
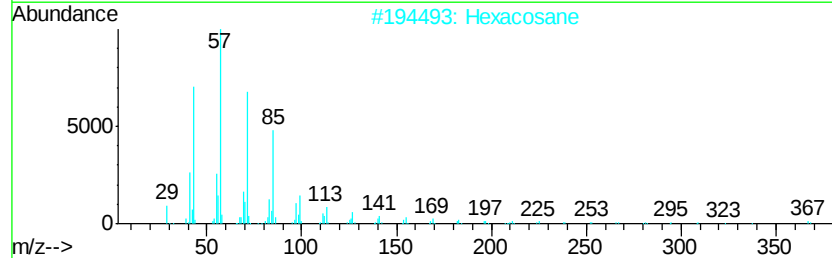
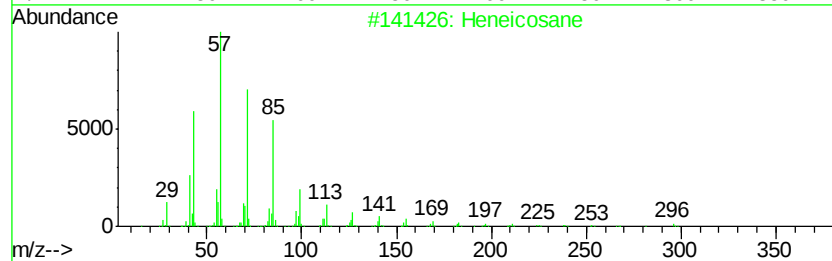
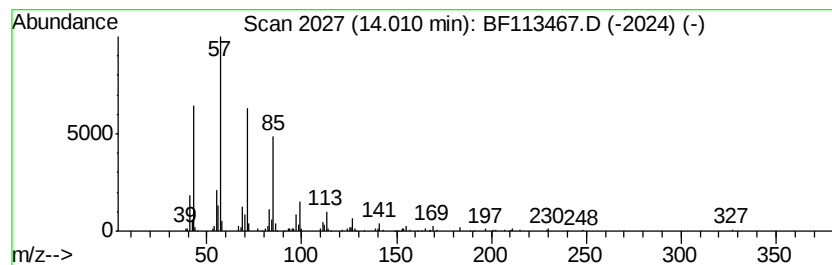
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	2.92 ng	166424	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Tetracosane		338	C24H50	000646-31-1	90
4	Octacosane		394	C28H58	000630-02-4	90
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90



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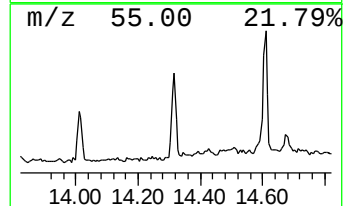
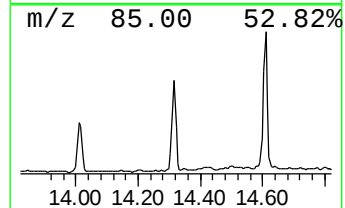
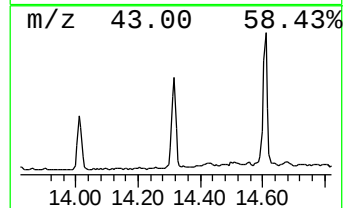
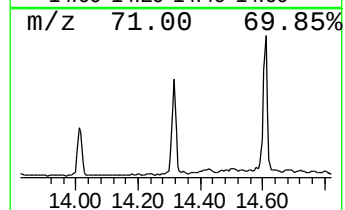
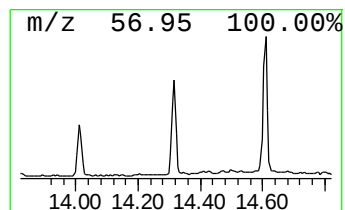
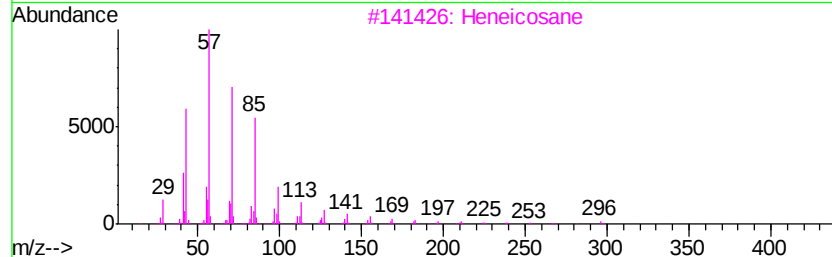
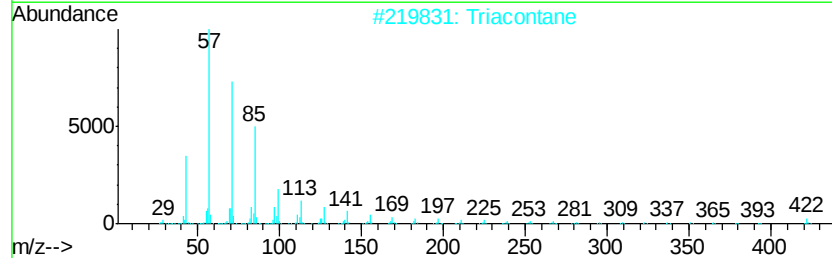
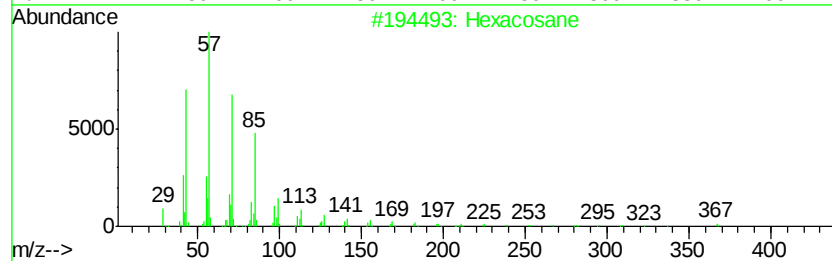
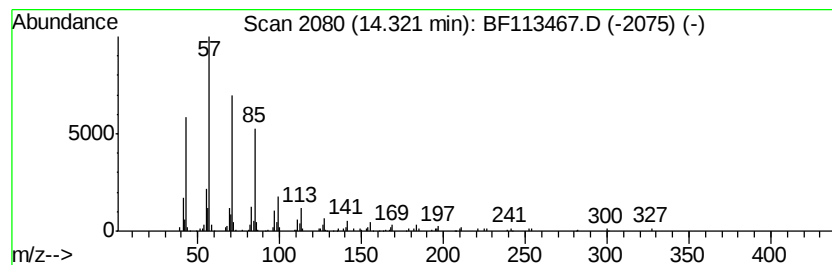
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.32	4.82 ng	274317	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



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GW-COMP

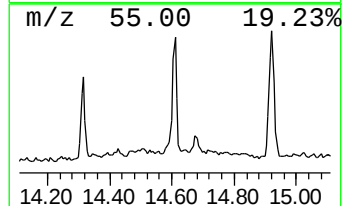
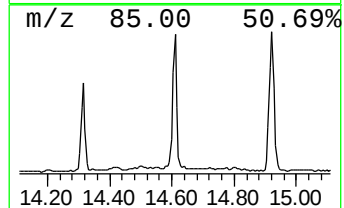
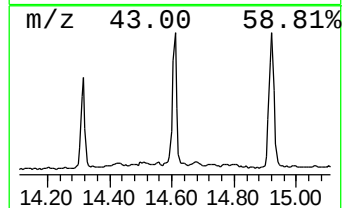
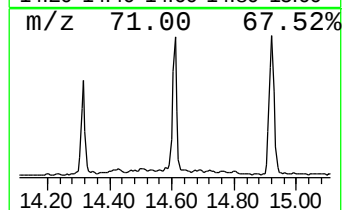
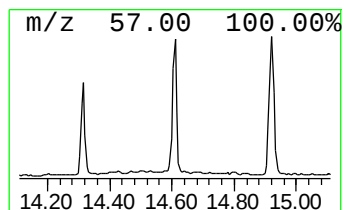
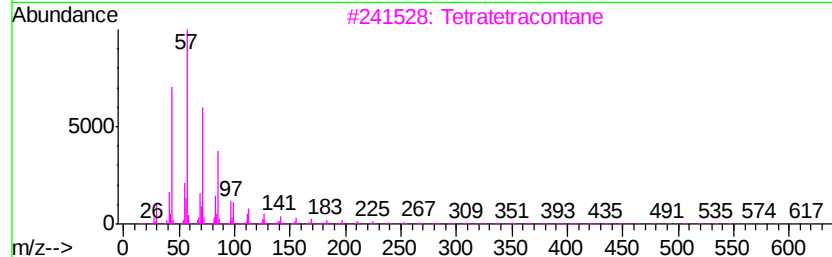
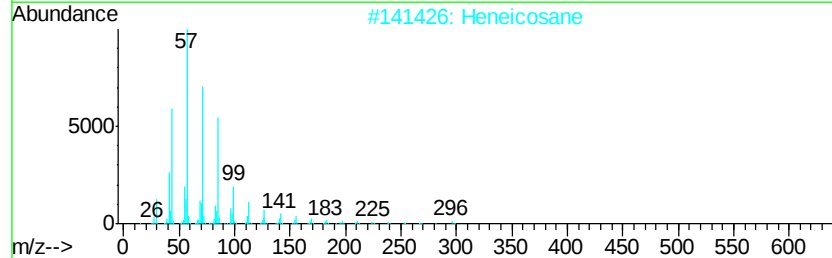
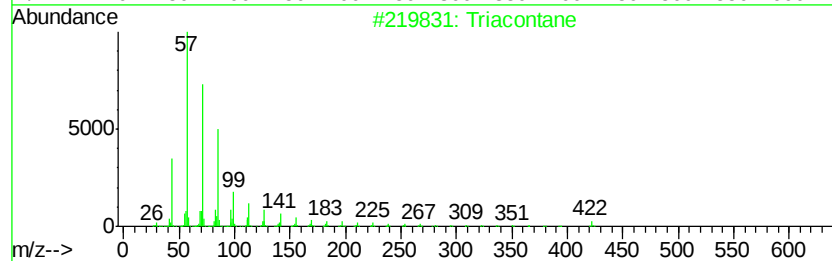
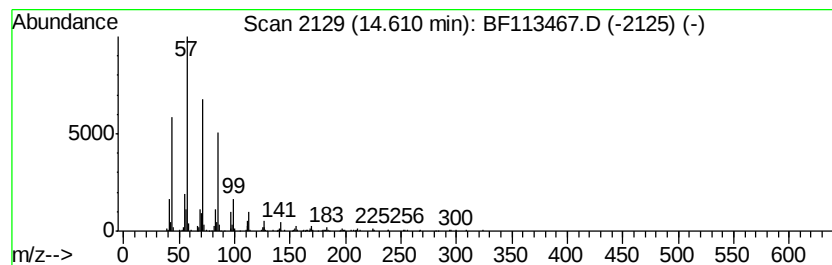
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Triacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	5.19 ng	427719	Perylene-d12	15.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			Heneicosane	296	C21H44	000629-94-7	90
3			Tetratetracontane	619	C44H90	007098-22-8	87
4			Heptacosane	380	C27H56	000593-49-7	83
5			Octacosane	394	C28H58	000630-02-4	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
Data File : BF113467.D
Acq On : 1 Apr 2019 15:32
Operator : JU/SJ
Sample : K2171-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GW-COMP

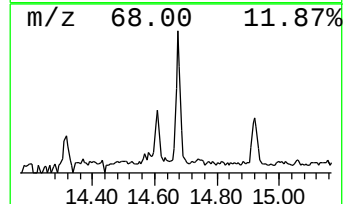
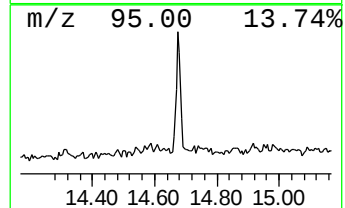
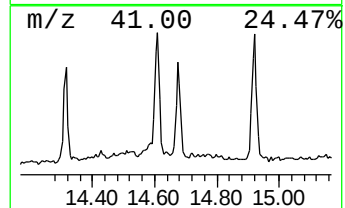
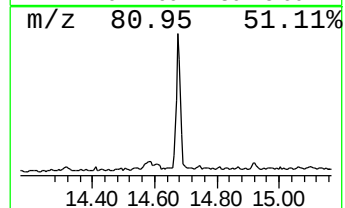
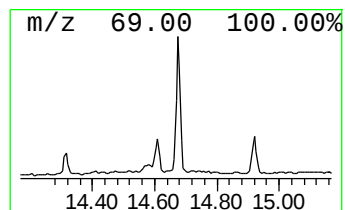
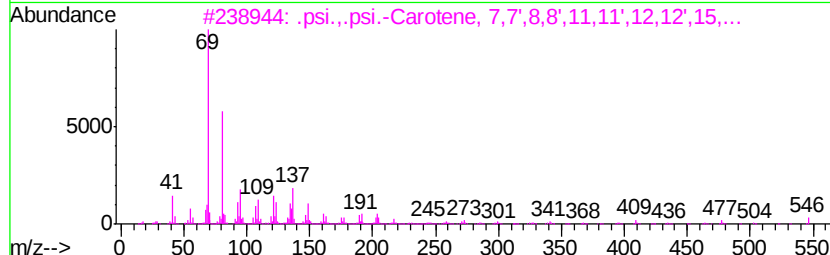
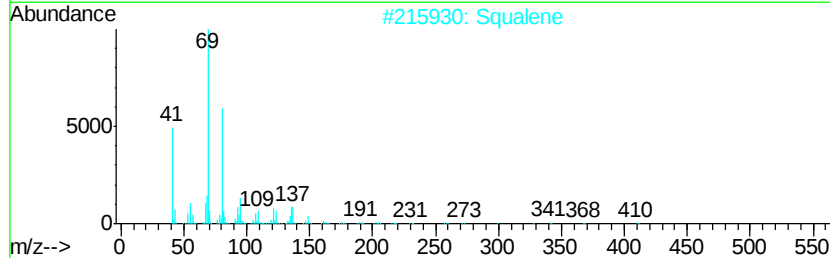
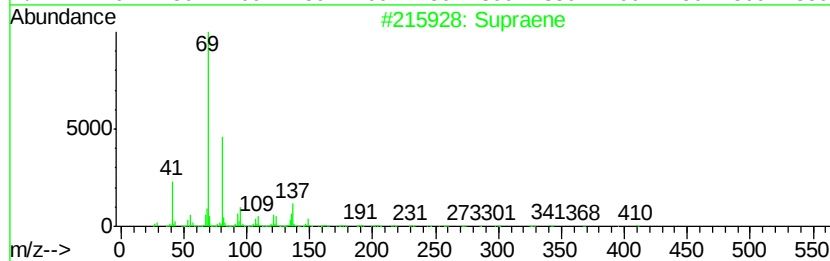
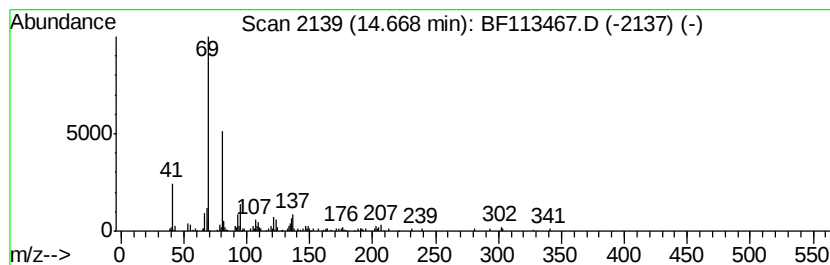
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Supraene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.45 ng	201654	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	91
3		.psi.,.psi.-Carotene, 7,7',8,8',...	547	C40H66	000502-62-5	83
4		trans-Farnesol	222	C15H26O	000106-28-5	64
5		1,5-Heptadiene, 2,6-dimethyl-	124	C9H16	006709-39-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
Data File : BF113467.D
Acq On : 1 Apr 2019 15:32
Operator : JU/SJ
Sample : K2171-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GW-COMP

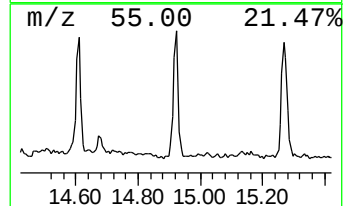
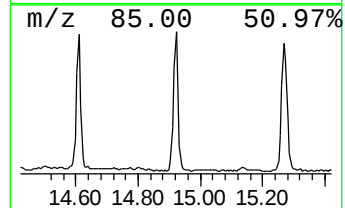
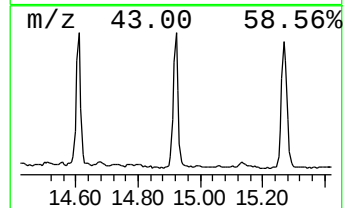
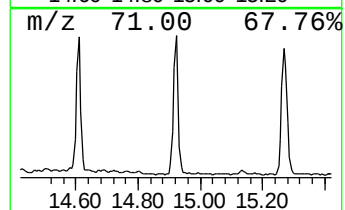
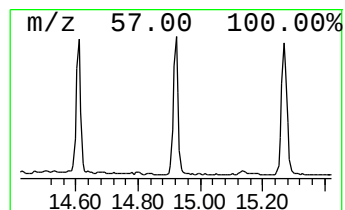
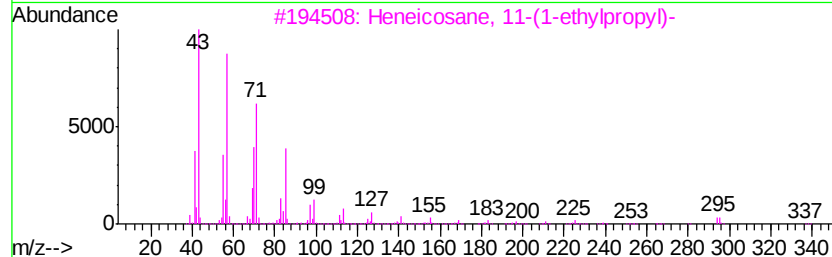
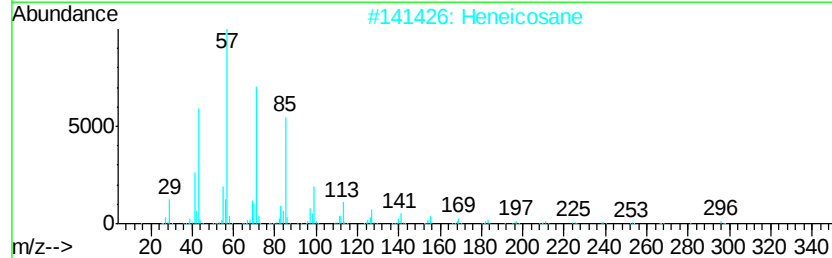
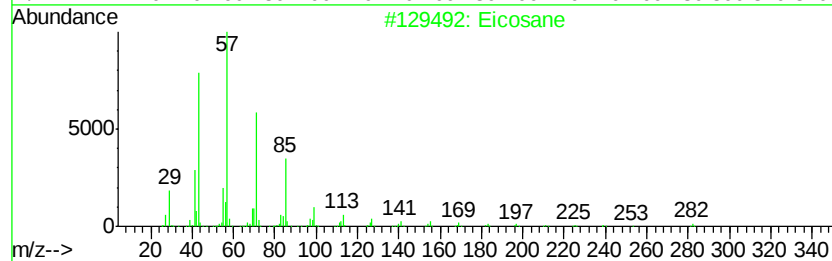
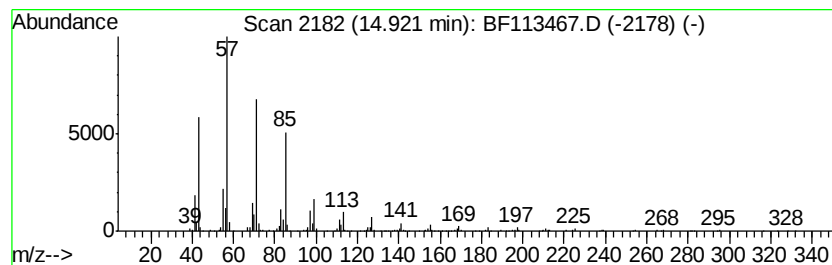
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	6.28 ng	516820	Perylene-d12	15.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	Heneicosane		296	C21H44	000629-94-7	91
3	Heneicosane, 11-(1-ethylpropyl)-		366	C26H54	055282-11-6	91
4	Triacontane		422	C30H62	000638-68-6	91
5	Hexadecane		226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
Data File : BF113467.D
Acq On : 1 Apr 2019 15:32
Operator : JU/SJ
Sample : K2171-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GW-COMP

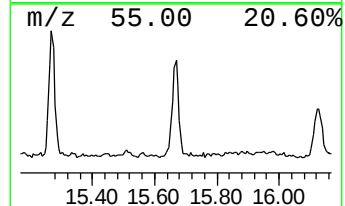
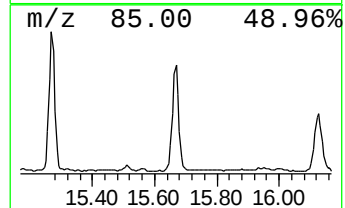
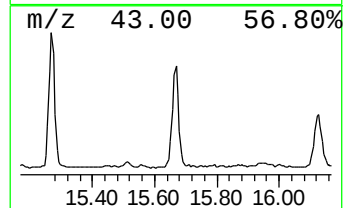
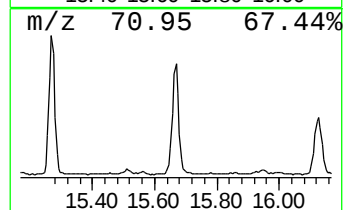
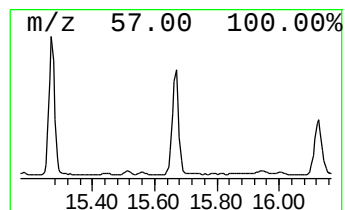
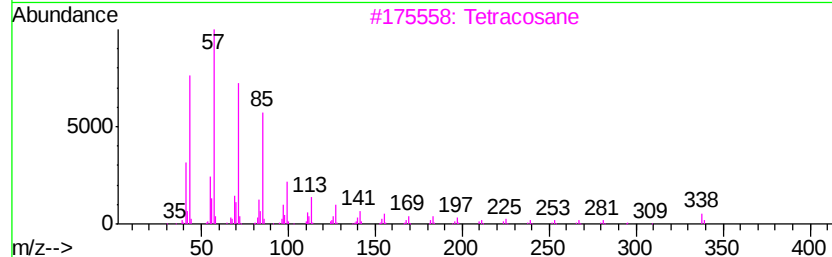
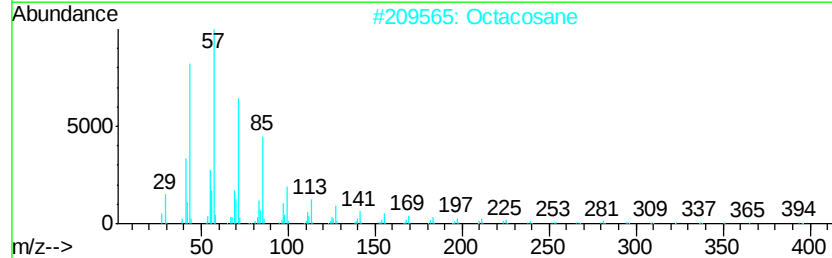
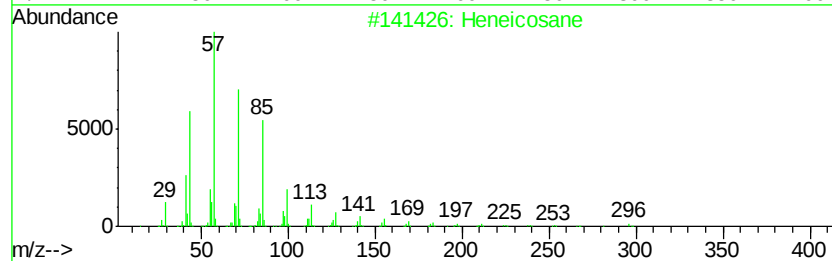
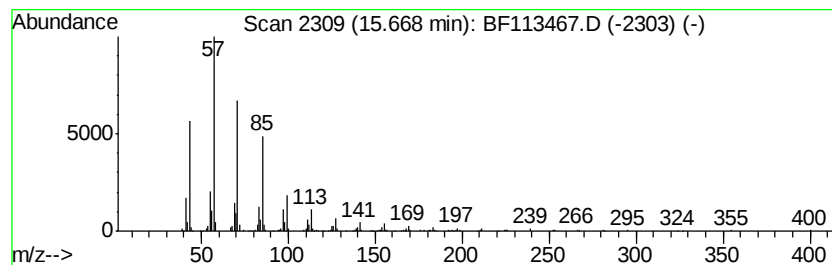
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-hexyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	6.26 ng	515611	Perylene-d12	15.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113467.D
 Acq On : 1 Apr 2019 15:32
 Operator : JU/SJ
 Sample : K2171-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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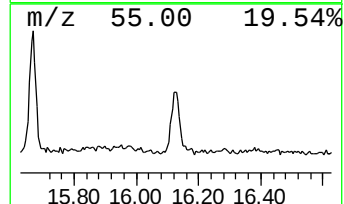
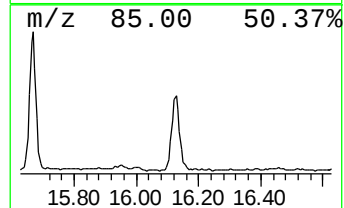
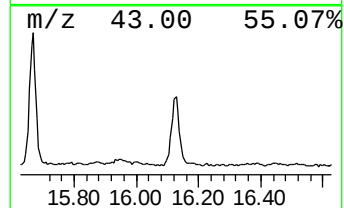
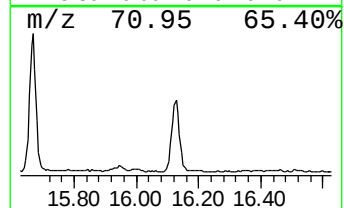
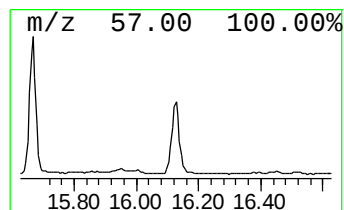
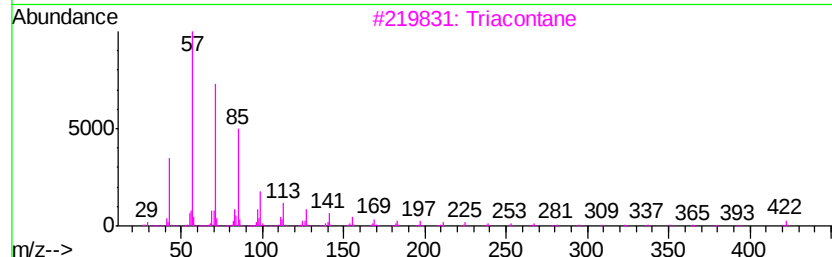
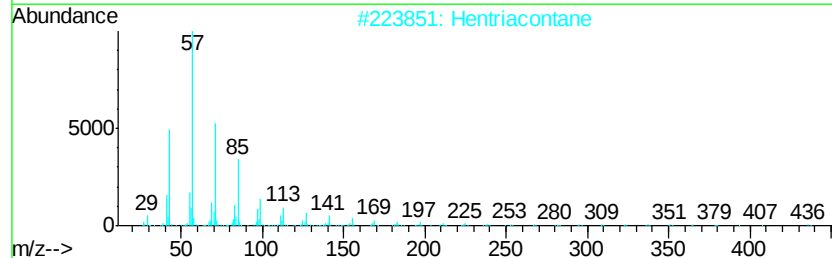
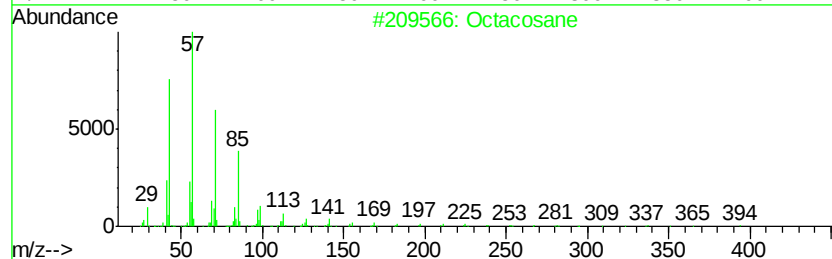
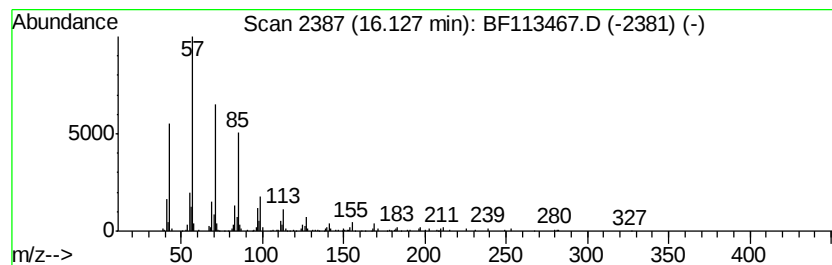
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	4.29 ng	353068	Perylene-d12	15.25

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			Triacontane	422	C30H62	000638-68-6	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113467.D
 Acq On : 1 Apr 2019 15:32
 Operator : JU/SJ
 Sample : K2171-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.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
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n-Hexadecanoic acid	11.75	3.4	ng	231251	4	11.22	1370750	20.0
Octadecanoic acid	12.53	5.9	ng	403997	4	11.22	1370750	20.0
Heneicosane	14.01	2.9	ng	166424	5	13.84	1139070	20.0
Hexacosane	14.32	4.8	ng	274317	5	13.84	1139070	20.0
Triacontane	14.61	5.2	ng	427719	6	15.25	1646780	20.0
Supraene	14.67	2.5	ng	201654	6	15.25	1646780	20.0
Eicosane	14.92	6.3	ng	516820	6	15.25	1646780	20.0
Heptadecane, 9-he...	15.67	6.3	ng	515611	6	15.25	1646780	20.0
Octacosane	16.13	4.3	ng	353068	6	15.25	1646780	20.0