

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012020\
 Data File : BF118596.D
 Acq On : 20 Jan 2020 17:36
 Operator : CG/JU
 Sample : L1181-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 93OAK-AVE-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-BF012020.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	2.216	14	22	29	rBV	2759833	3868464	28.85%	2.765%
2	5.104	505	513	522	rBV	1636982	2313794	17.26%	1.654%
3	5.504	575	581	584	rBV	9535613	10327603	77.02%	7.382%
4	6.504	745	751	755	rBV	9324209	11341539	84.59%	8.107%
5	6.651	771	776	779	rBV	10993528	11144043	83.11%	7.966%
6	6.881	811	815	818	rBV	3328260	2781993	20.75%	1.989%
7	7.040	837	842	845	rVB	9508143	8685727	64.78%	6.208%
8	7.440	904	910	913	rBV	6534177	6935171	51.72%	4.957%
9	8.163	1028	1033	1036	rBV	3584672	3543581	26.43%	2.533%
10	9.239	1210	1216	1219	rBV	13248835	12890432	96.14%	9.214%
11	9.628	1278	1282	1285	rBV	1049109	834105	6.22%	0.596%
12	9.916	1326	1331	1335	rBV	5029158	4416208	32.94%	3.157%
13	10.704	1459	1465	1468	rBV	9733955	9488530	70.77%	6.782%
14	11.404	1579	1584	1587	rBV	4961003	4577829	34.14%	3.272%
15	11.469	1590	1595	1598	rVB2	168219	157229	1.17%	0.112%
16	11.927	1667	1673	1689	rBV	880708	971656	7.25%	0.695%
17	12.710	1802	1806	1819	rBV	265583	349745	2.61%	0.250%
18	12.992	1849	1854	1857	rBV	12994269	13408171	100.00%	9.584%
19	13.563	1948	1951	1958	rVB	207656	221701	1.65%	0.158%
20	13.880	2002	2005	2020	rBV2	2001979	2403633	17.93%	1.718%
21	14.033	2027	2031	2035	rBV	3386491	3300239	24.61%	2.359%
22	14.210	2056	2061	2070	rBV	1807638	1772323	13.22%	1.267%
23	14.516	2107	2113	2123	rBV	3445635	3378412	25.20%	2.415%
24	14.827	2160	2166	2176	rBV	4134820	4482518	33.43%	3.204%
25	15.168	2218	2224	2235	rBV	3562317	4583908	34.19%	3.276%
26	15.504	2276	2281	2285	rBV	2154771	2793203	20.83%	1.997%
27	15.557	2285	2290	2300	rVB	2888773	4059109	30.27%	2.901%
28	15.998	2358	2365	2370	rBV	1786476	2953222	22.03%	2.111%
29	16.510	2445	2452	2470	rVB	774045	1554308	11.59%	1.111%
30	17.115	2549	2555	2561	rBV2	181446	365455	2.73%	0.261%

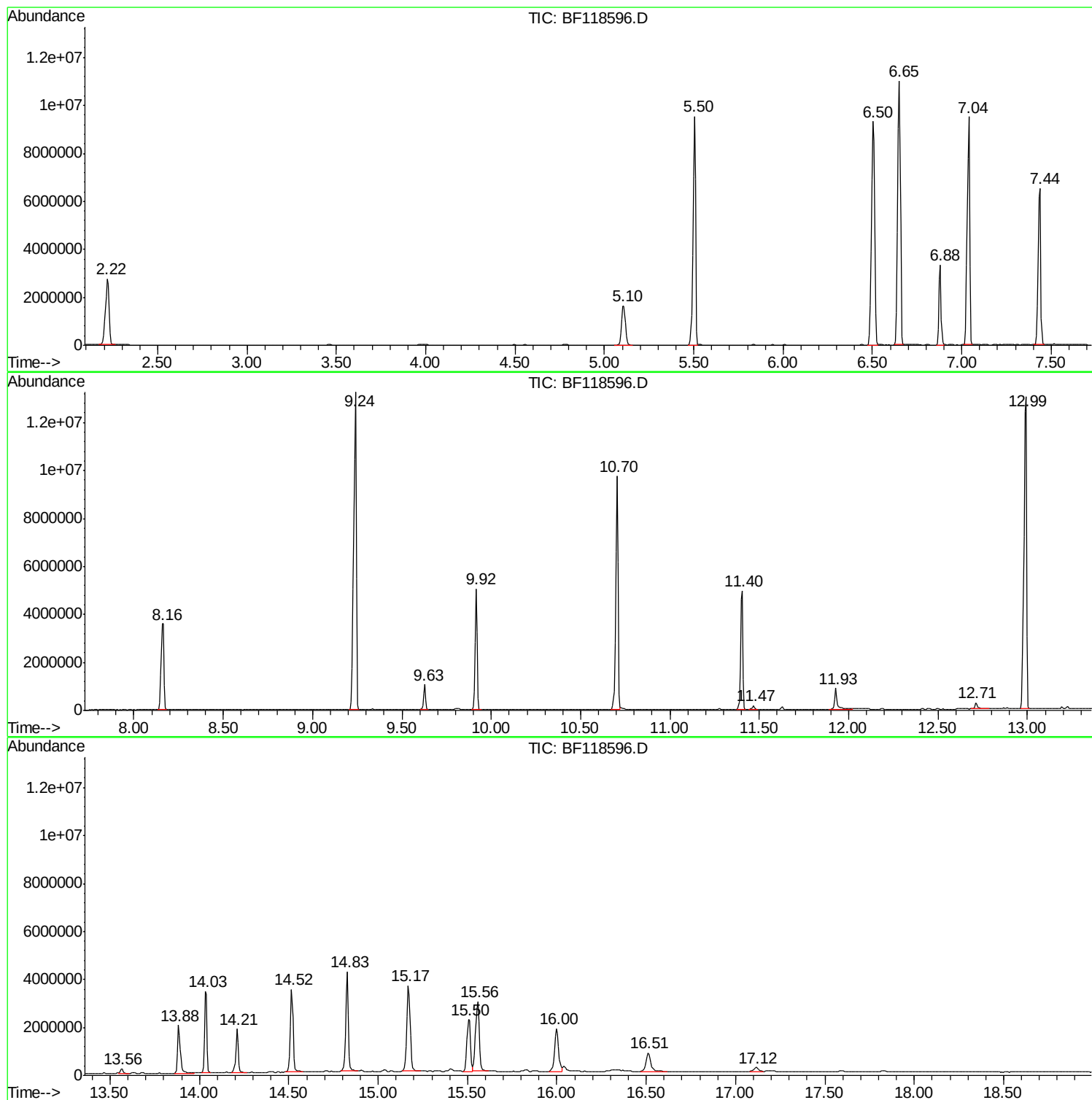
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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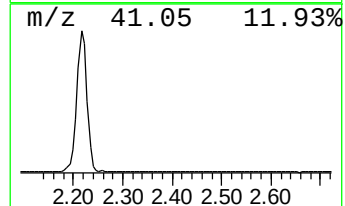
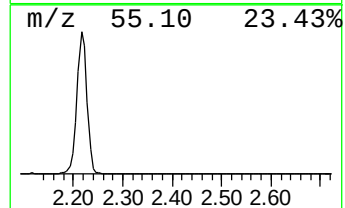
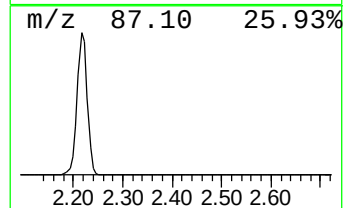
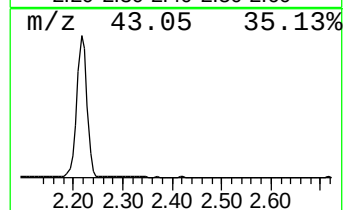
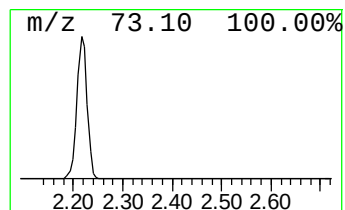
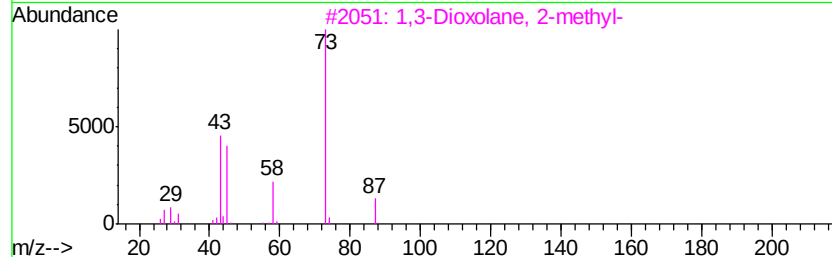
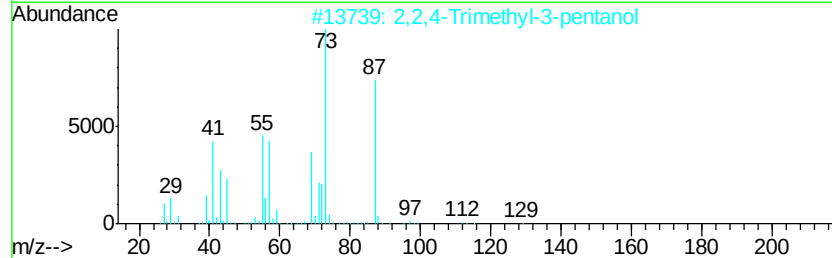
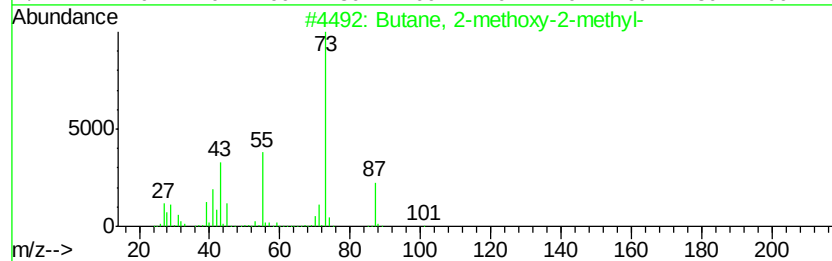
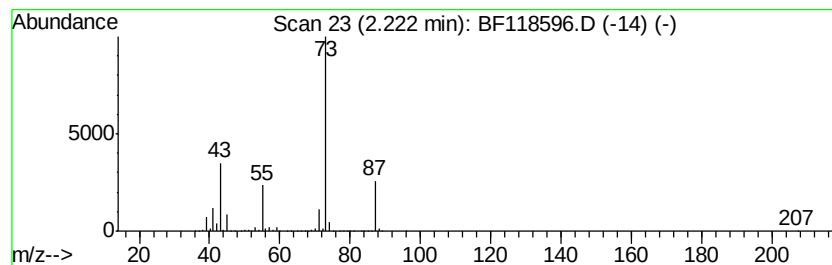
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.22	27.81 ng	3868460	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22



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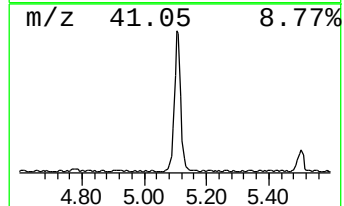
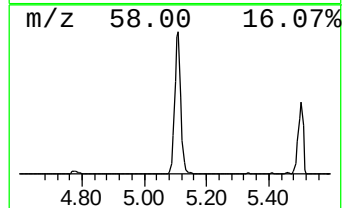
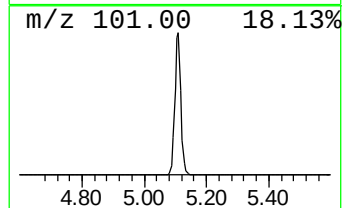
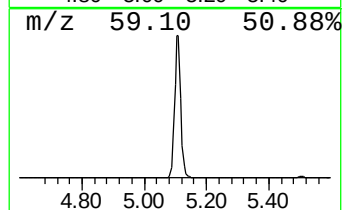
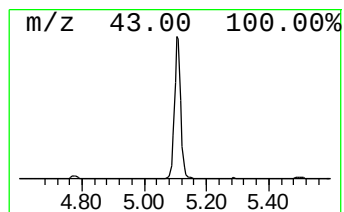
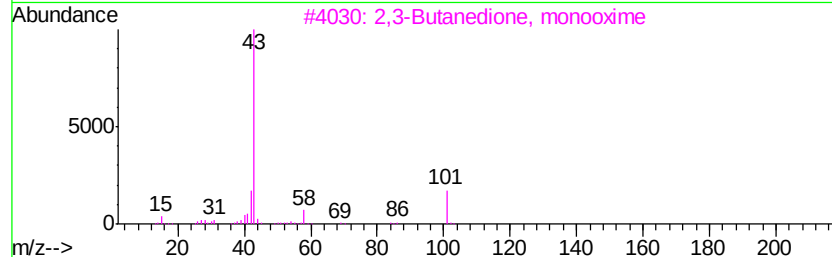
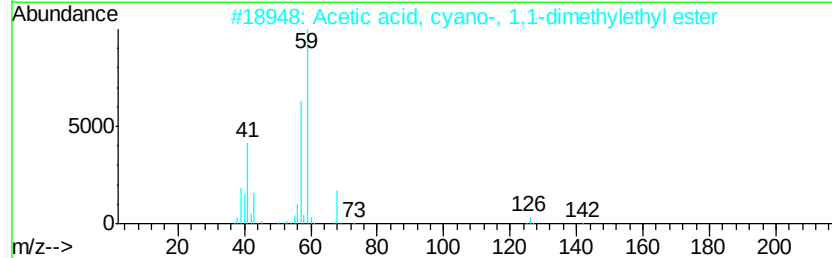
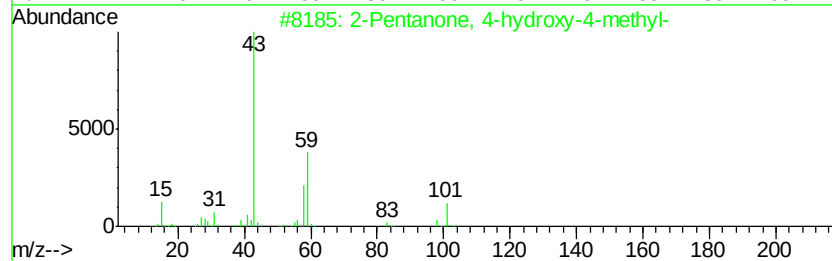
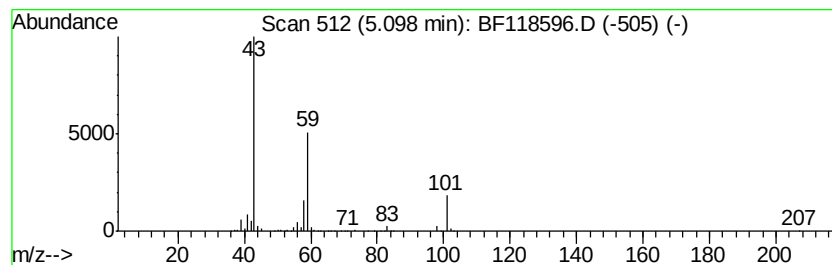
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	16.63 ng	2313790	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4		Acetone	58	C3H6O	000067-64-1	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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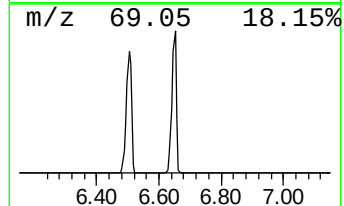
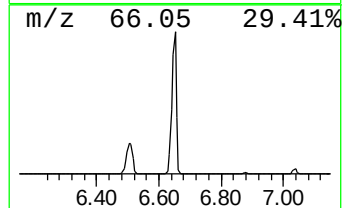
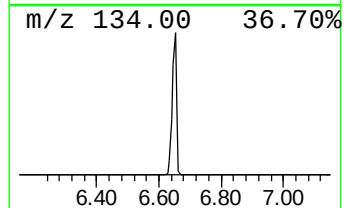
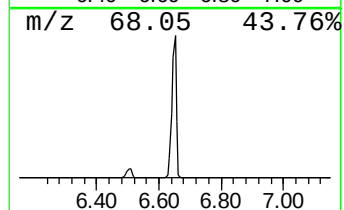
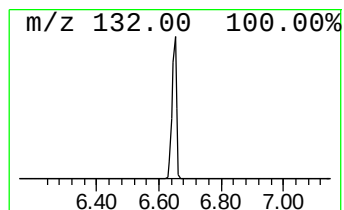
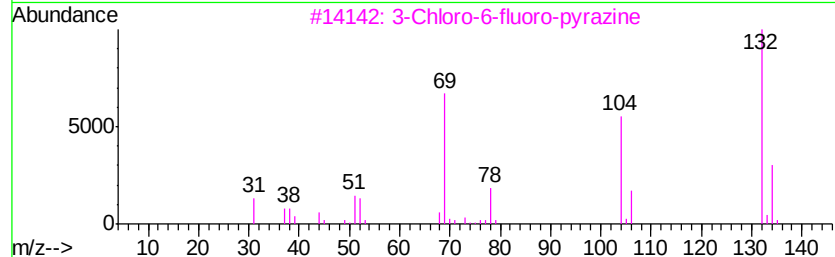
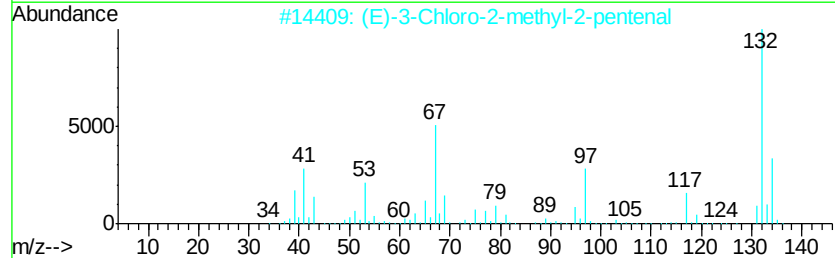
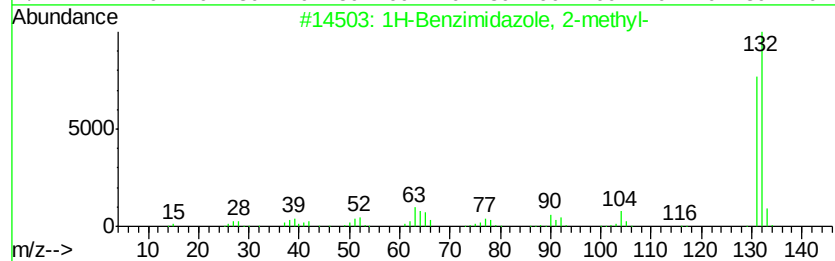
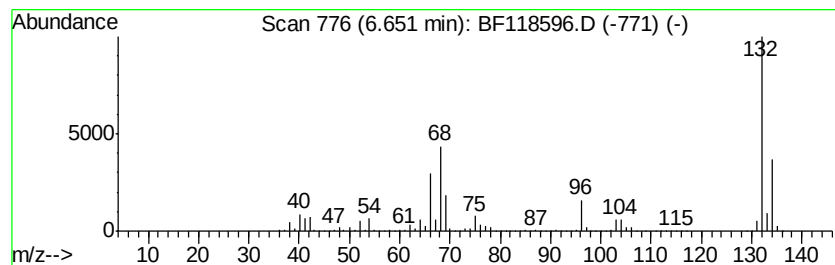
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	80.12 ng	11144000	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		1-Isouquinolinemethanol, .alpha.-...	191	C12H17NO	114608-92-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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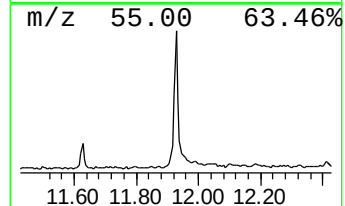
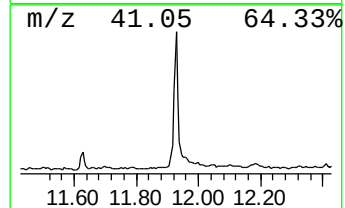
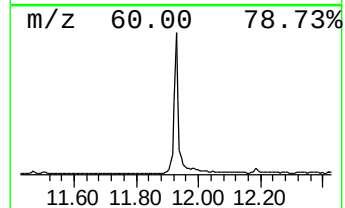
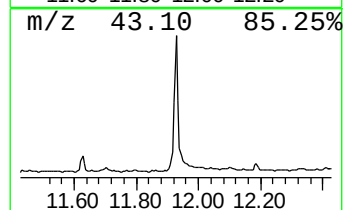
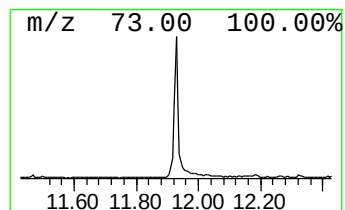
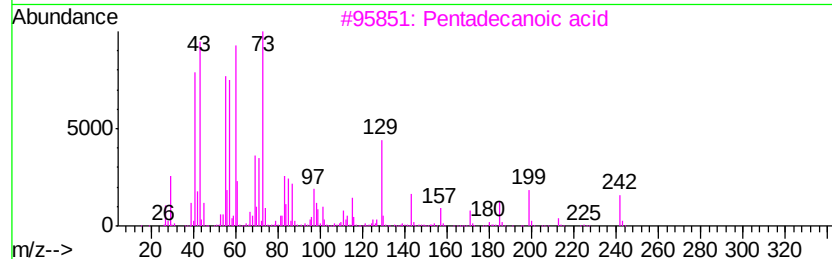
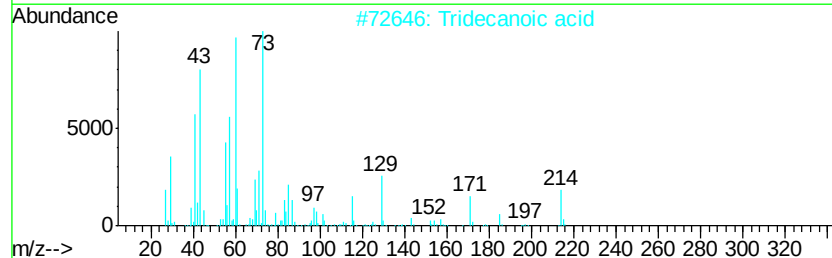
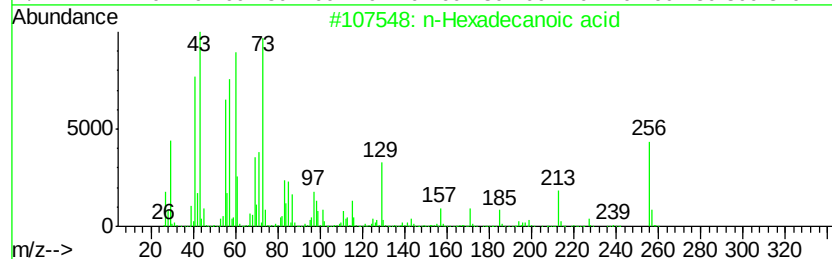
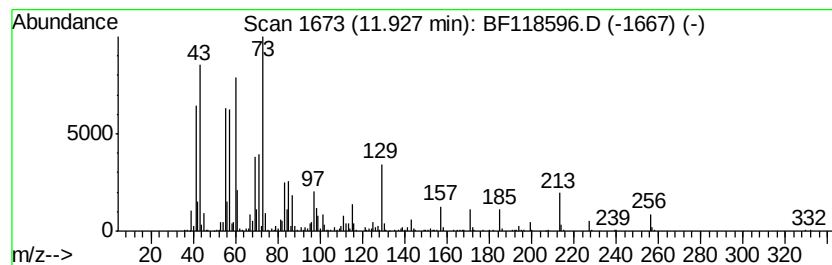
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	4.25 ng	971656	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Octadecanoic acid	284	C18H36O2	000057-11-4	87
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



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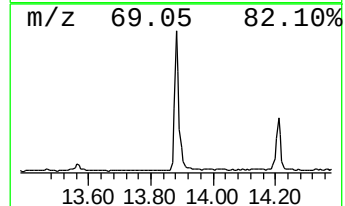
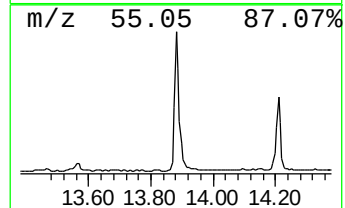
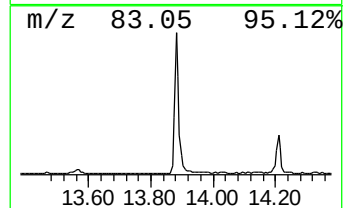
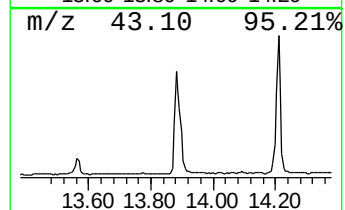
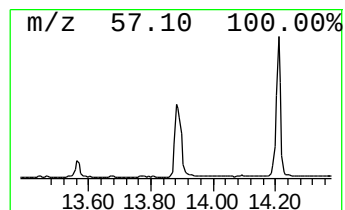
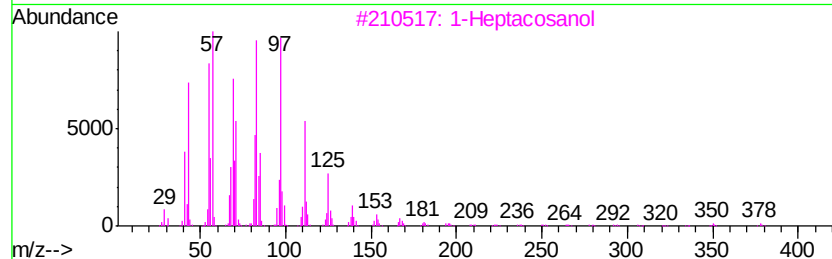
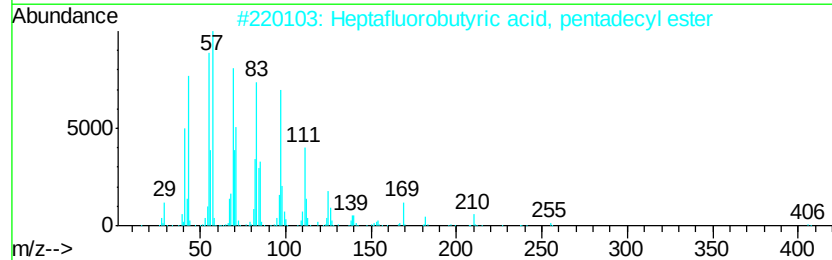
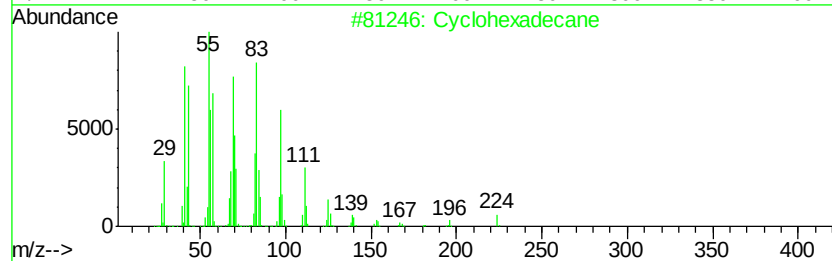
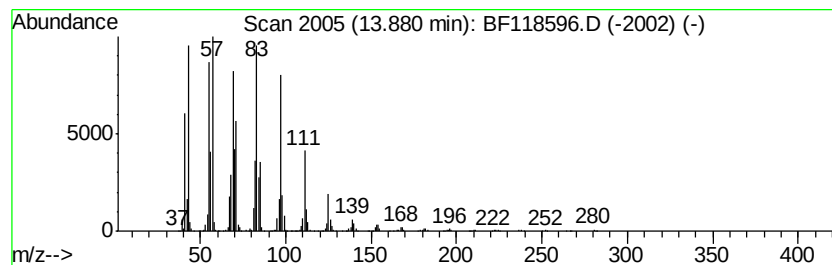
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 Peak Number 6 Cyclohexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	14.57 ng	2403630	Chrysene-d12	14.03
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 94
2		Heptafluorobutyric acid, pentade...	424 C19H31F7O2	959261-23-5 94
3		1-Heptacosanol	396 C27H56O	002004-39-9 94
4		Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1 94
5		Dichloroacetic acid, heptadecyl ...	366 C19H36Cl2O2	1000282-98-2 94



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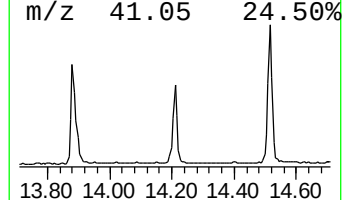
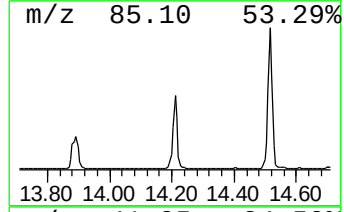
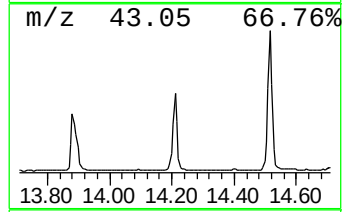
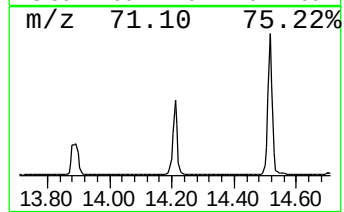
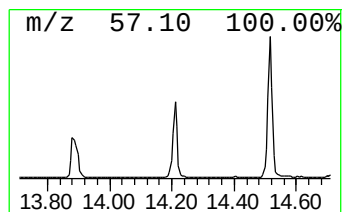
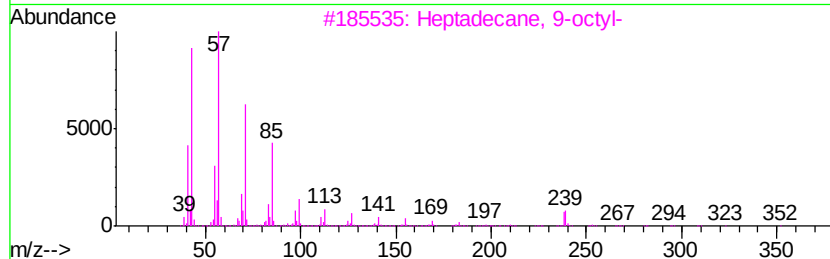
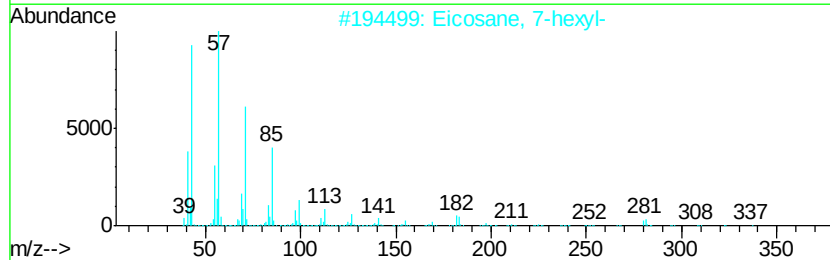
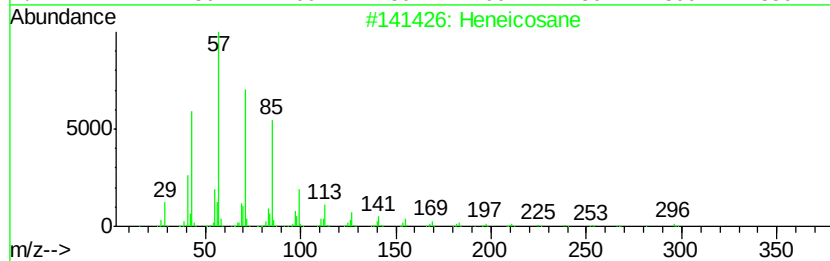
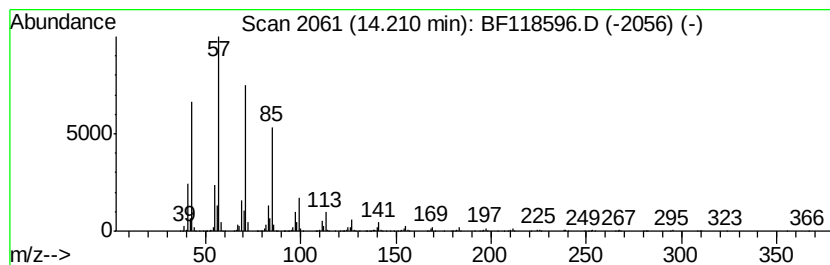
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	10.74 ng	1772320	Chrysene-d12	14.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Hexatriacontane		507	C36H74	000630-06-8	91
5	Octadecane		254	C18H38	000593-45-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118596.D
Acq On : 20 Jan 2020 17:36
Operator : CG/JU
Sample : L1181-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

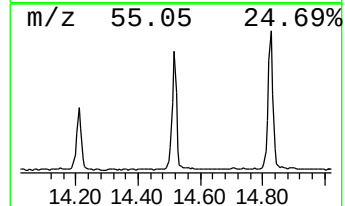
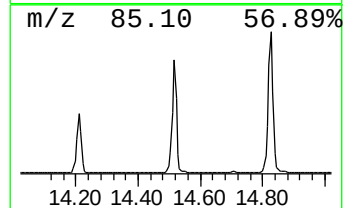
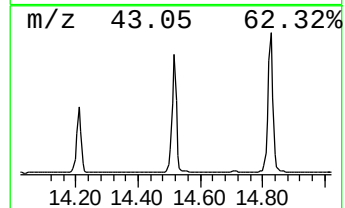
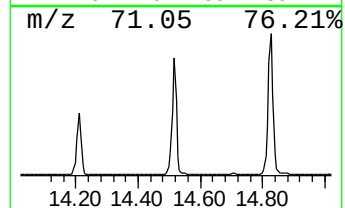
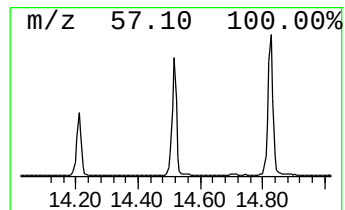
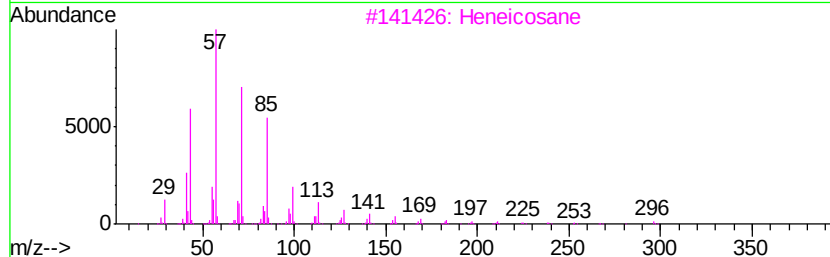
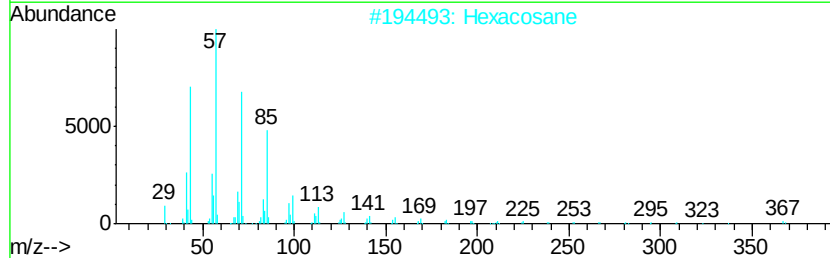
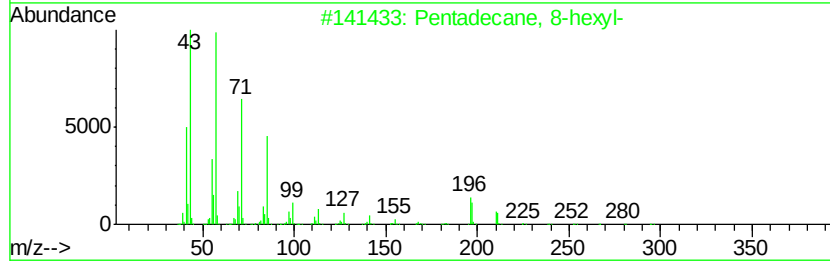
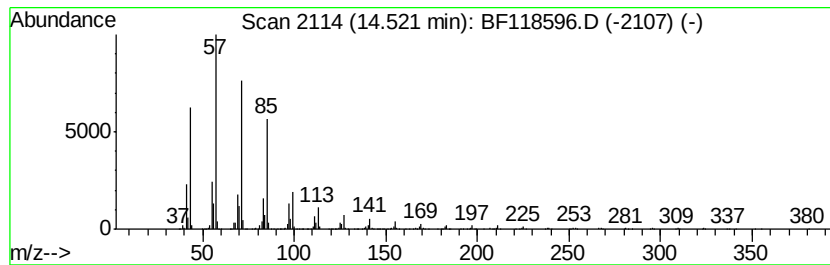
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 8-hexyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	20.47 ng	3378410	Chrysene-d12	14.03
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pentadecane, 8-hexyl-	296 C21H44	013475-75-7 94
2		Hexacosane	366 C26H54	000630-01-3 91
3		Heneicosane	296 C21H44	000629-94-7 91
4		Octacosane	394 C28H58	000630-02-4 91
5		Heptadecane, 9-octyl-	352 C25H52	007225-64-1 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
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Sample : L1181-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

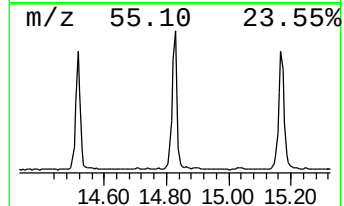
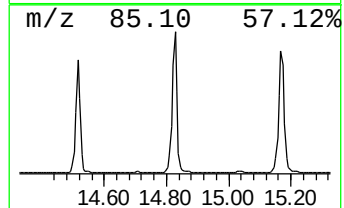
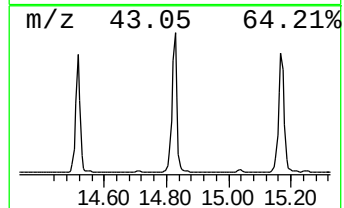
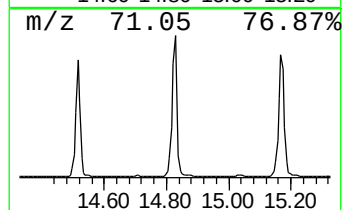
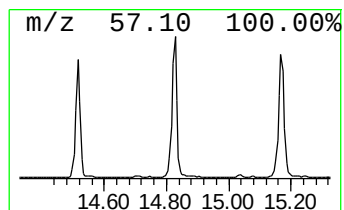
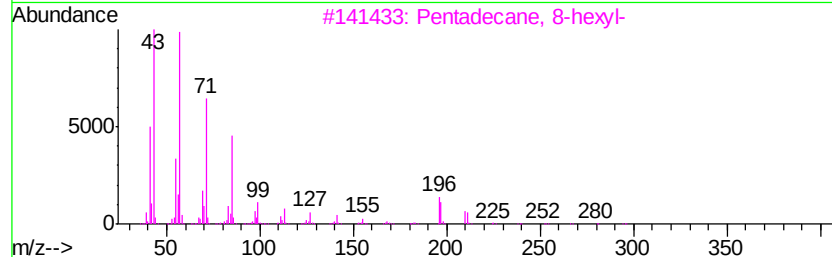
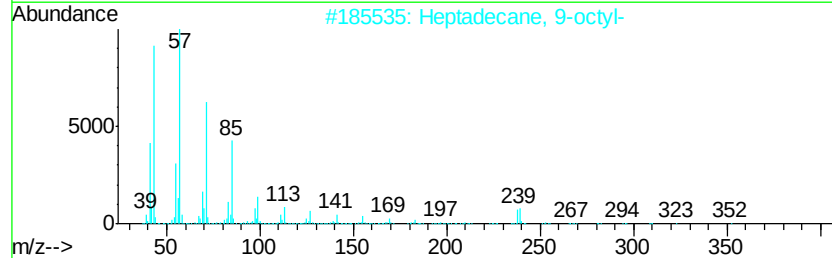
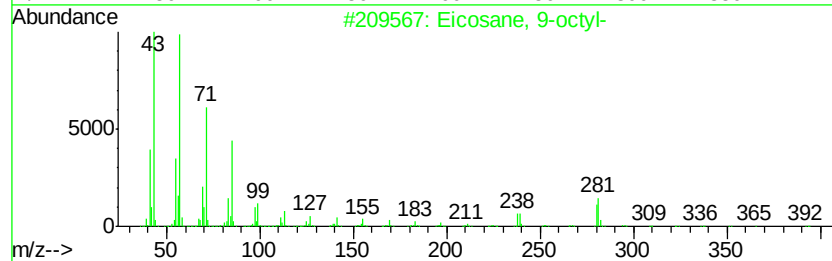
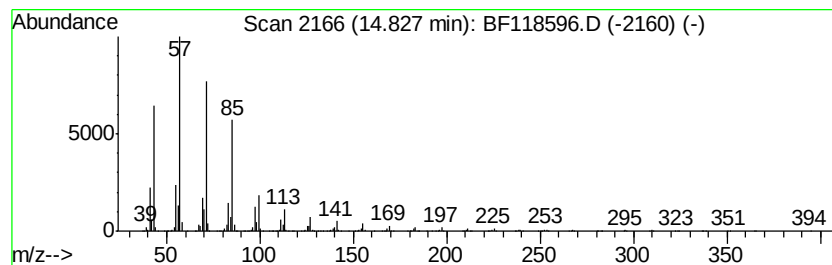
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane, 9-octyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	32.10 ng	4482520	Perylene-d12	15.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane, 9-octyl-	394 C28H58	013475-77-9	94
2	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
3	Pentadecane, 8-hexyl-	296 C21H44	013475-75-7	91
4	2-methyloctacosane	408 C29H60	1000376-72-8	91
5	Hexacosane	366 C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118596.D
Acq On : 20 Jan 2020 17:36
Operator : CG/JU
Sample : L1181-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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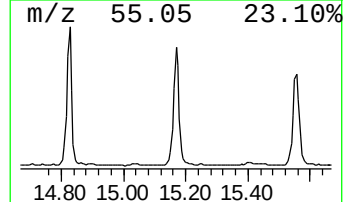
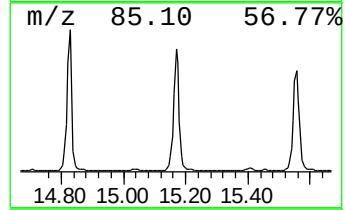
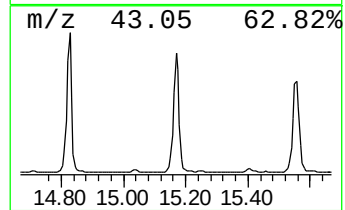
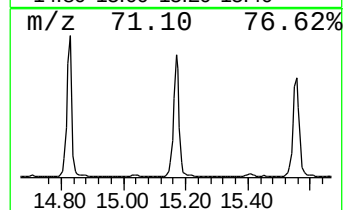
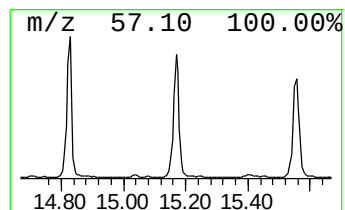
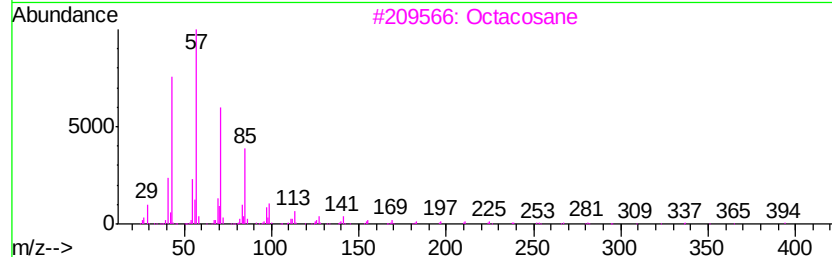
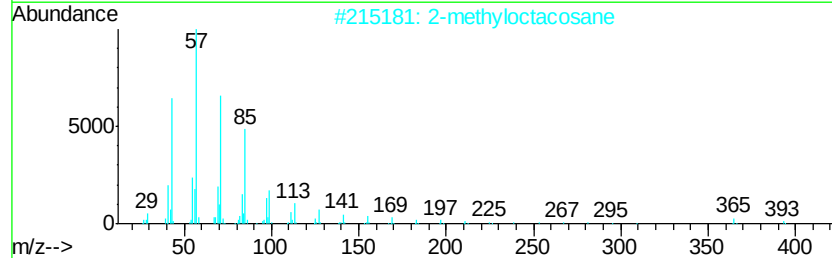
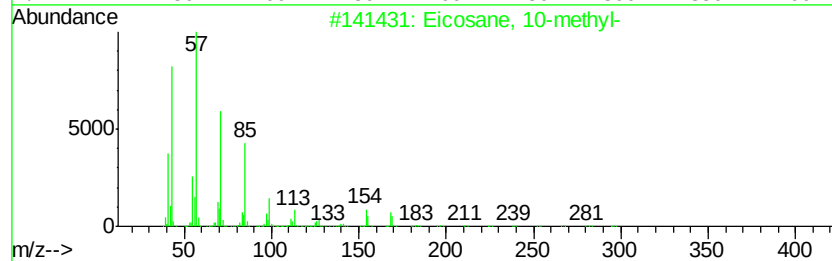
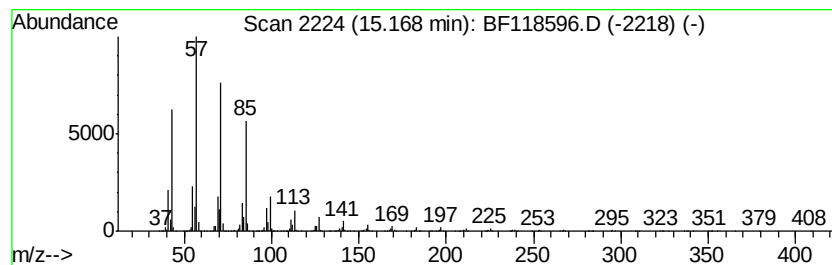
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane, 10-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	32.82 ng	4583910	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118596.D
Acq On : 20 Jan 2020 17:36
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Sample : L1181-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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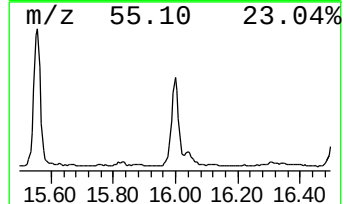
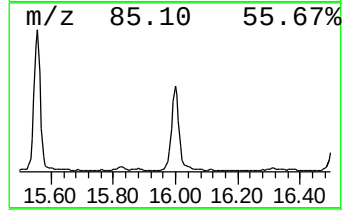
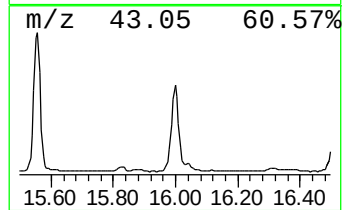
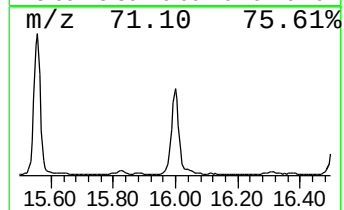
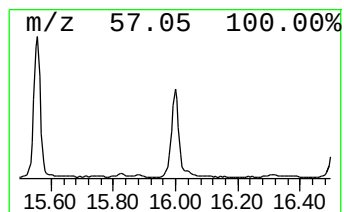
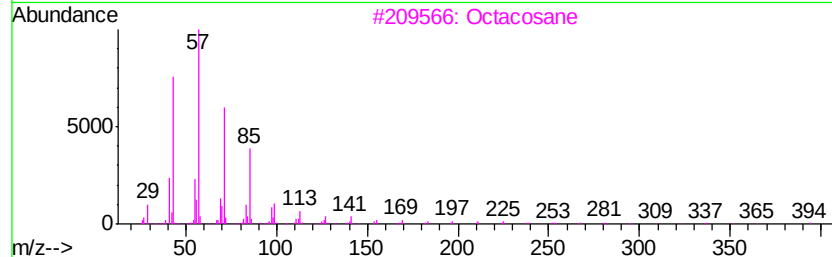
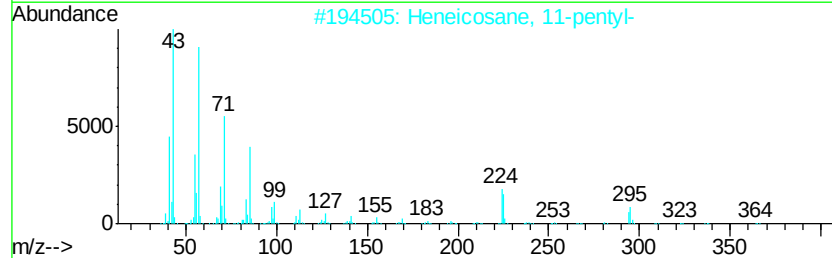
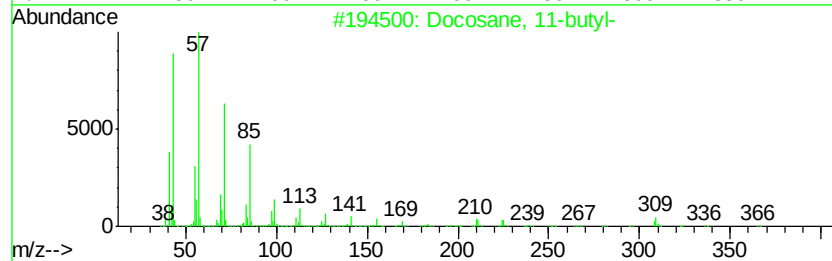
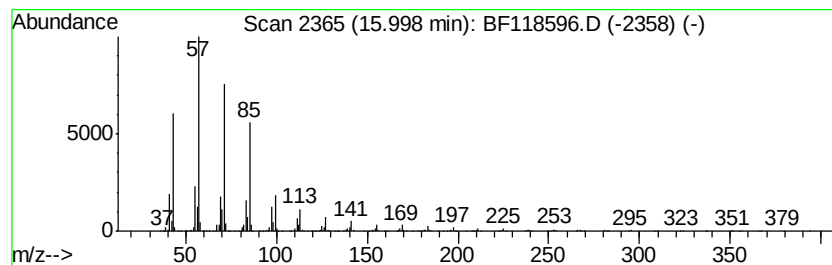
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Docosane, 11-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	21.15 ng	2953220	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
3		Octacosane	394	C28H58	000630-02-4	91
4		Docosane, 9-butyl-	366	C26H54	055282-14-9	91
5		Docosane, 5-butyl-	366	C26H54	055282-16-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118596.D
Acq On : 20 Jan 2020 17:36
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Misc :
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Instrument :
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ClientSampleId :
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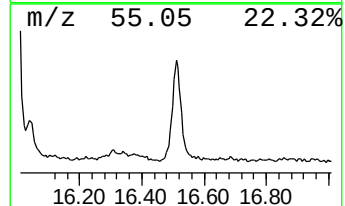
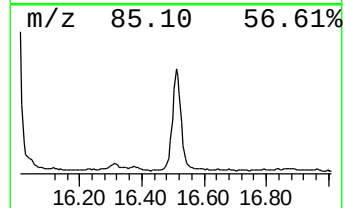
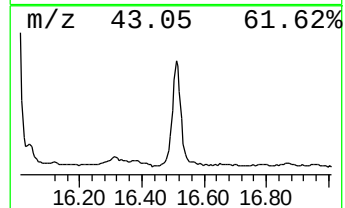
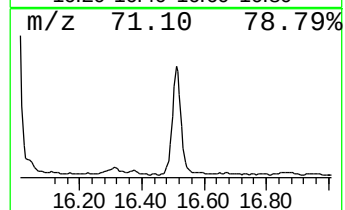
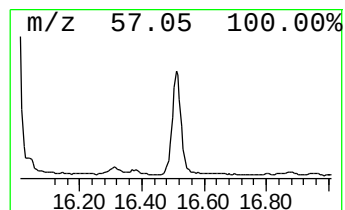
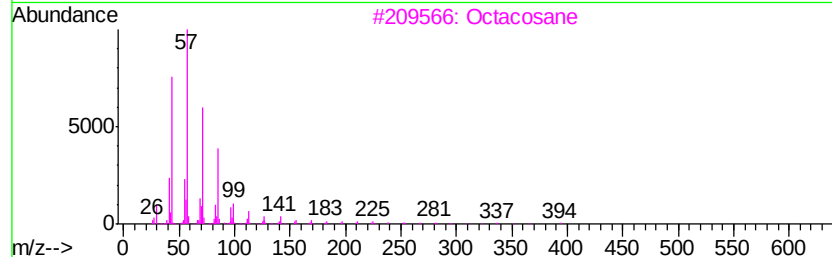
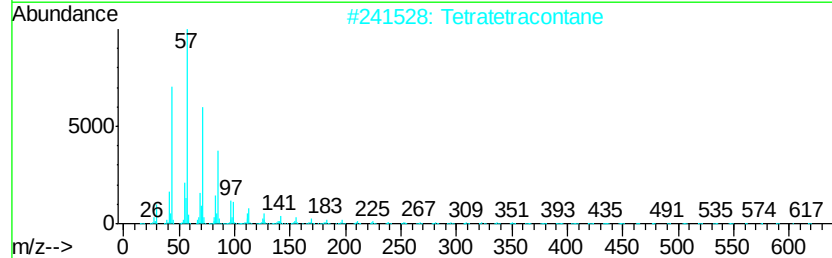
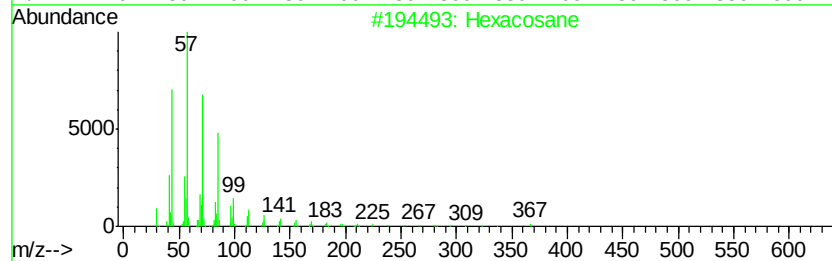
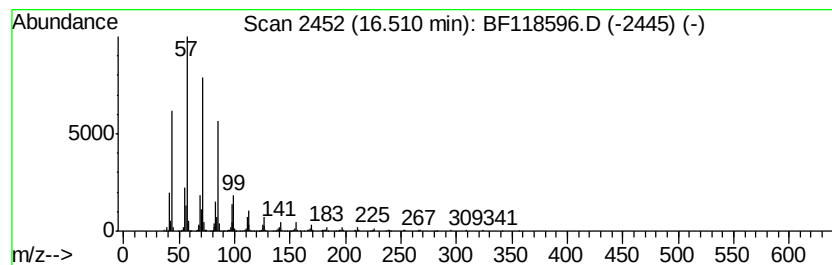
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	11.13 ng	1554310	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118596.D
Acq On : 20 Jan 2020 17:36
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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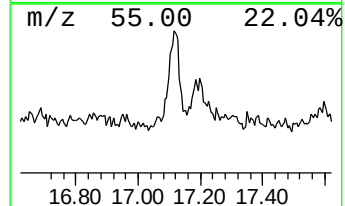
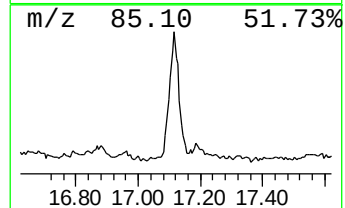
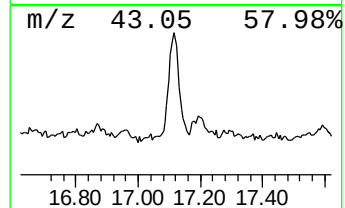
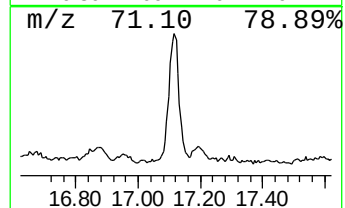
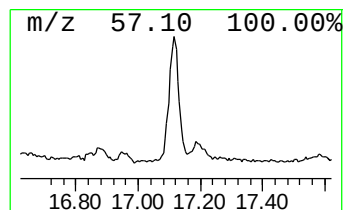
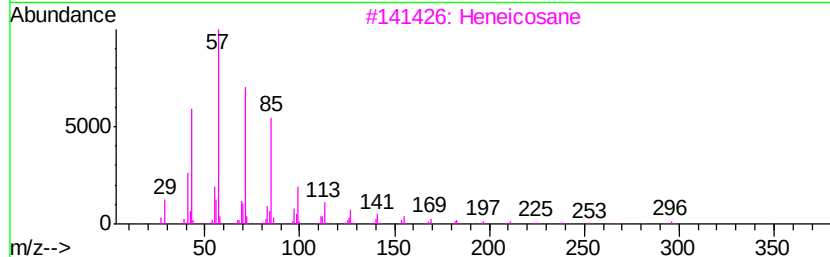
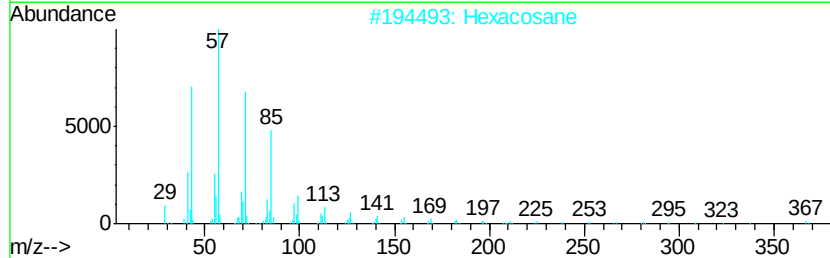
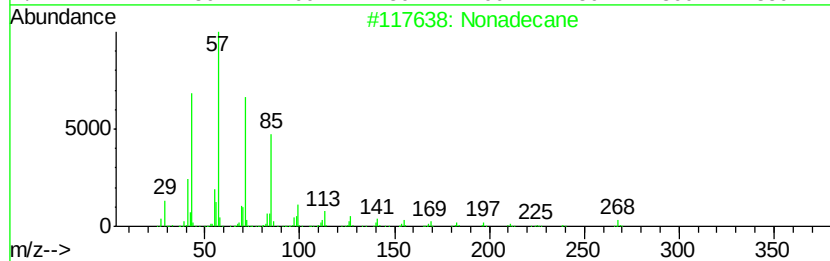
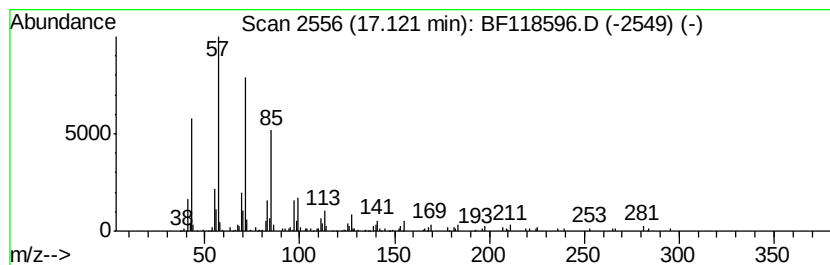
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.12	2.62 ng	365455	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Hexacosane	366	C26H54	000630-01-3	94
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012020\
 Data File : BF118596.D
 Acq On : 20 Jan 2020 17:36
 Operator : CG/JU
 Sample : L1181-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012020.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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2-Pentanone, 4-hy...	5.10	16.6	ng	2313790	1	6.88	2781990	20.0
unknown6.65	6.65	80.1	ng	11144000	1	6.88	2781990	20.0
n-Hexadecanoic acid	11.93	4.3	ng	971656	4	11.40	4577830	20.0
Cyclohexadecane	13.88	14.6	ng	2403630	5	14.03	3300240	20.0
Heneicosane	14.21	10.7	ng	1772320	5	14.03	3300240	20.0
Pentadecane, 8-he...	14.52	20.5	ng	3378410	5	14.03	3300240	20.0
Eicosane, 9-octyl-	14.83	32.1	ng	4482520	6	15.51	2793200	20.0
Eicosane, 10-methyl-	15.17	32.8	ng	4583910	6	15.51	2793200	20.0
Docosane, 11-butyl-	16.00	21.1	ng	2953220	6	15.51	2793200	20.0
Hexacosane	16.51	11.1	ng	1554310	6	15.51	2793200	20.0
Nonadecane	17.12	2.6	ng	365455	6	15.51	2793200	20.0