

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
 Data File : BF118603.D
 Acq On : 20 Jan 2020 20:51
 Operator : CG/JU
 Sample : L1154-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-0-15

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	2.222	14	23	30	rBV	2677716	3700415	28.73%	3.041%
2	5.104	506	513	521	rVB	1640592	2260404	17.55%	1.858%
3	5.504	575	581	584	rBV	9585098	10269739	79.74%	8.441%
4	6.504	745	751	755	rBV	9411425	11456995	88.96%	9.417%
5	6.651	770	776	779	rBV	10887406	11230068	87.20%	9.230%
6	6.881	811	815	818	rBV	3330671	2894394	22.47%	2.379%
7	7.040	837	842	845	rVB	9434734	8697523	67.53%	7.149%
8	7.440	904	910	913	rBV	6525941	6996257	54.32%	5.750%
9	8.163	1028	1033	1036	rBV	3971976	3760824	29.20%	3.091%
10	9.239	1210	1216	1219	rBV	12995168	12879138	100.00%	10.586%
11	9.628	1277	1282	1285	rBV	1069517	901281	7.00%	0.741%
12	9.916	1326	1331	1335	rBV	5325137	4562538	35.43%	3.750%
13	10.704	1459	1465	1468	rBV	9369700	8885978	69.00%	7.303%
14	11.398	1579	1583	1587	rBV	4829312	4478001	34.77%	3.681%
15	11.469	1589	1595	1598	rVB3	192234	176310	1.37%	0.145%
16	11.927	1669	1673	1678	rBV	822905	743386	5.77%	0.611%
17	12.710	1803	1806	1818	rBV2	266613	320560	2.49%	0.263%
18	12.986	1848	1853	1856	rBV	12419072	12139942	94.26%	9.978%
19	13.880	2002	2005	2018	rBV	1897609	1844617	14.32%	1.516%
20	14.033	2027	2031	2035	rBV	3315245	3001525	23.31%	2.467%
21	14.210	2056	2061	2073	rBV	501063	483761	3.76%	0.398%
22	14.516	2108	2113	2125	rBV	911221	973446	7.56%	0.800%
23	14.821	2160	2165	2175	rBV	1152268	1275828	9.91%	1.049%
24	15.168	2218	2224	2234	rBV	1161803	1473656	11.44%	1.211%
25	15.504	2276	2281	2285	rBV	2146568	2747825	21.34%	2.258%
26	15.551	2285	2289	2297	rVB	965842	1362647	10.58%	1.120%
27	15.998	2358	2365	2369	rBV	700160	1104191	8.57%	0.908%
28	16.039	2369	2372	2382	rVB3	178187	297484	2.31%	0.245%
29	16.510	2446	2452	2460	rBV	313750	567658	4.41%	0.467%
30	17.115	2549	2555	2562	rBV2	89648	181109	1.41%	0.149%

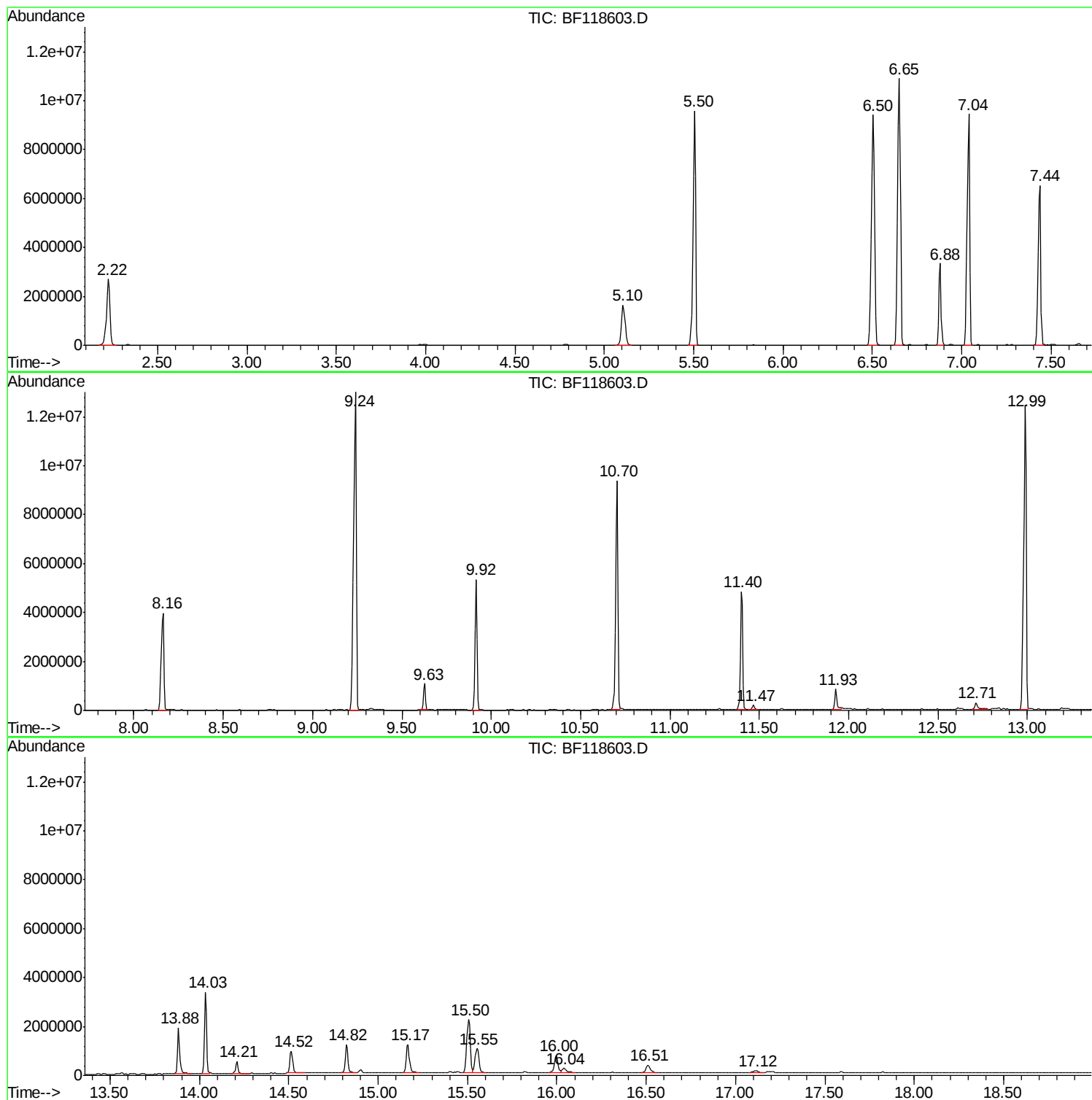
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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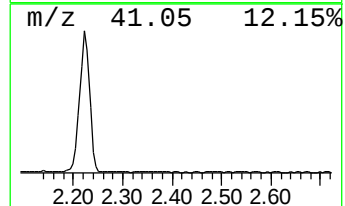
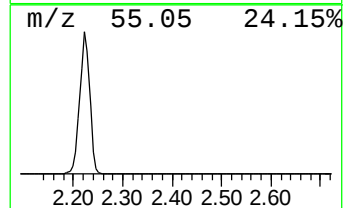
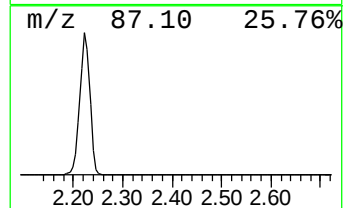
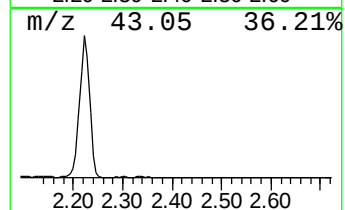
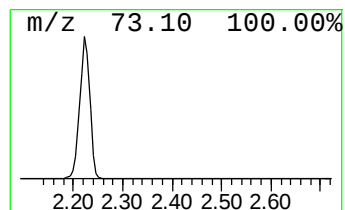
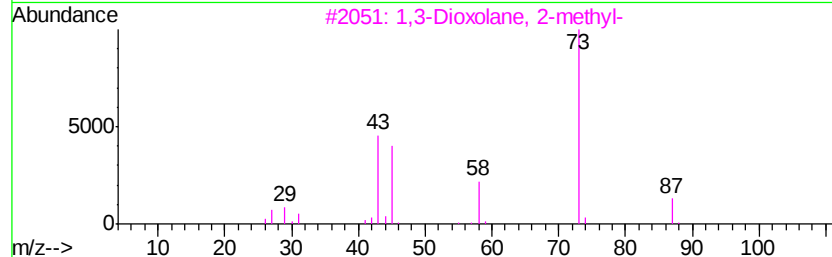
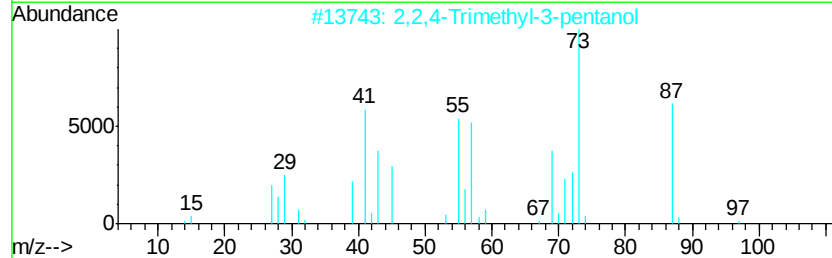
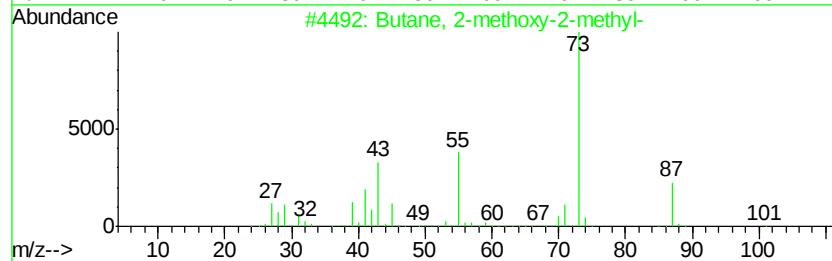
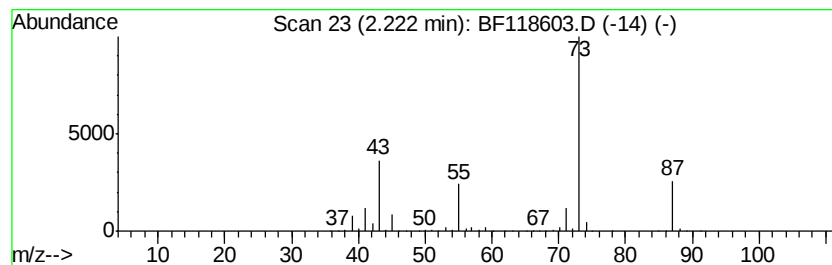
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.22	25.57 ng	3700420	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	56
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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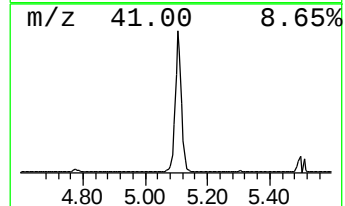
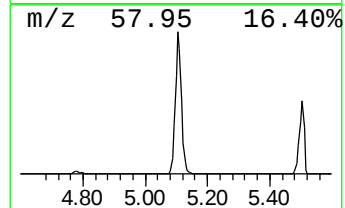
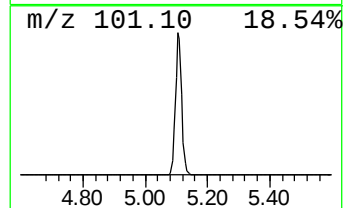
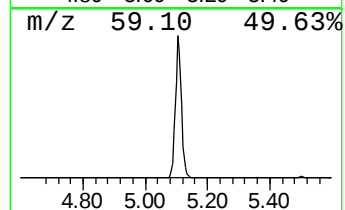
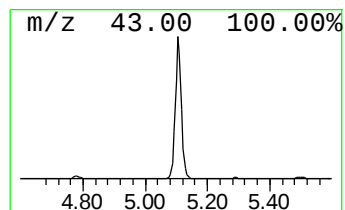
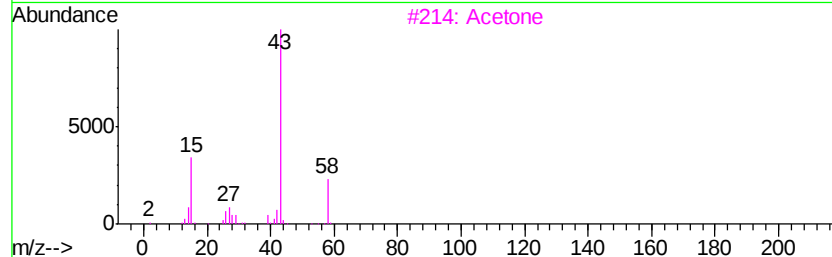
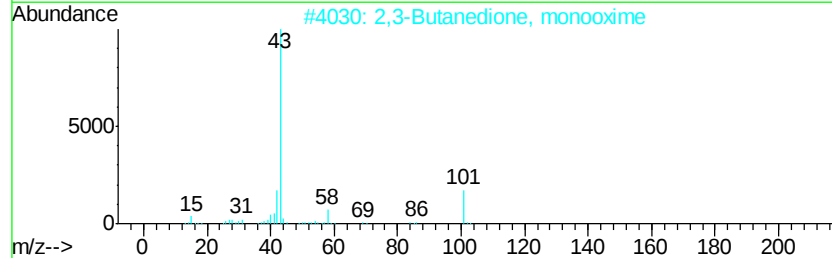
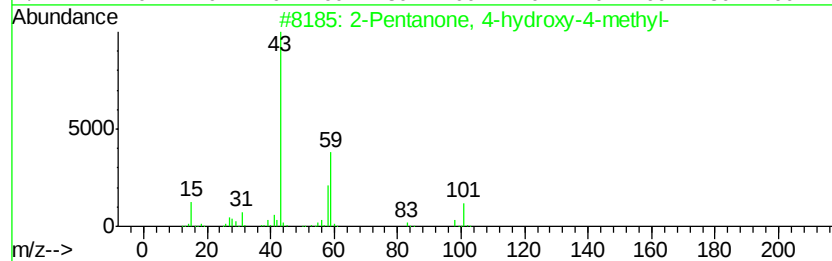
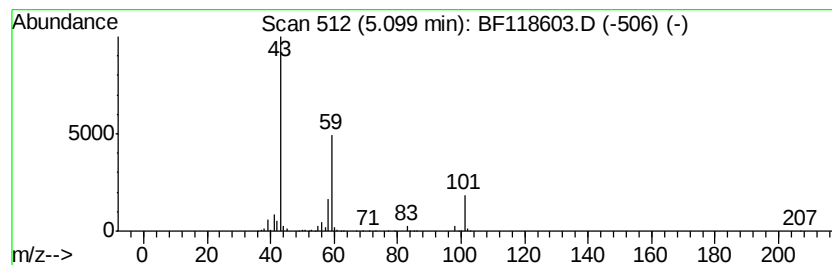
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	15.62 ng	2260400	1,4-Dichlorobenzene-d4	6.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	59
2	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	16
3	Acetone	58	C3H6O		000067-64-1	12
4	1-Propen-2-ol, acetate	100	C5H8O2		000108-22-5	10
5	Morpholine, 4-methyl-	101	C5H11NO		000109-02-4	9



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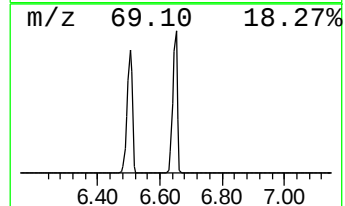
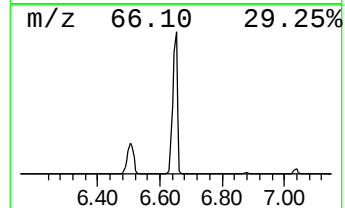
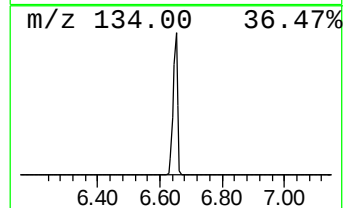
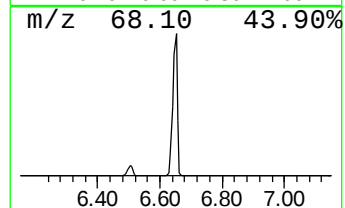
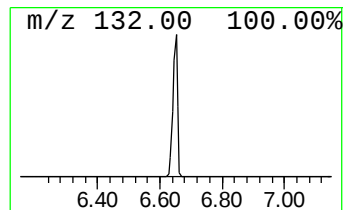
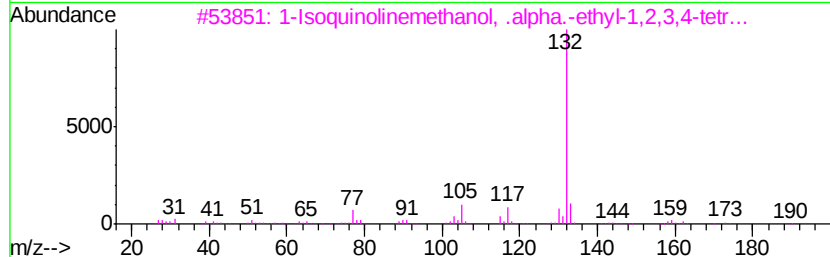
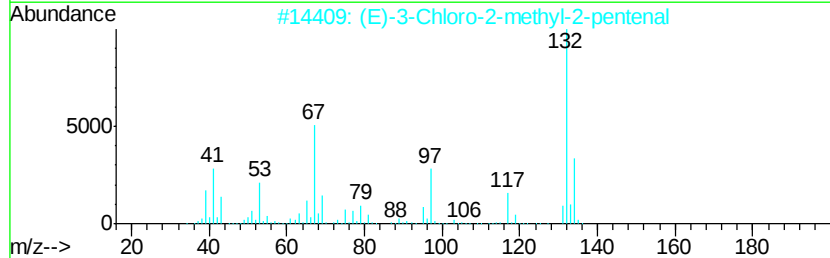
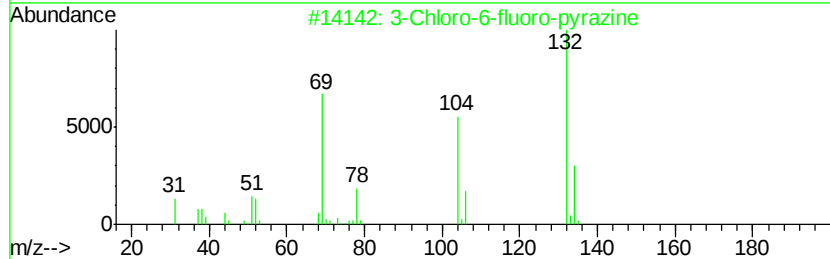
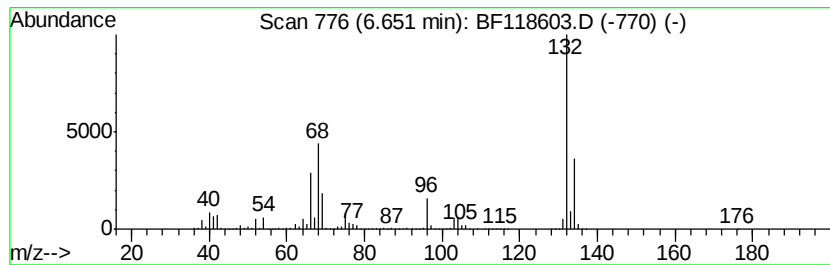
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	77.60 ng	11230100	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		1-Isoquinolinemethanol, .alpha.-...	191	C12H17NO	114608-92-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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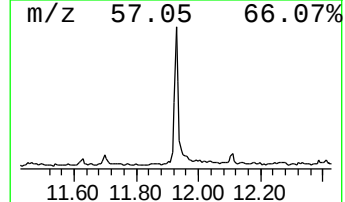
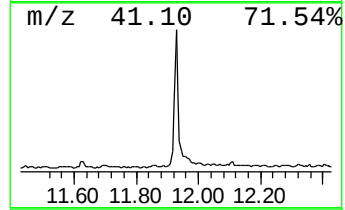
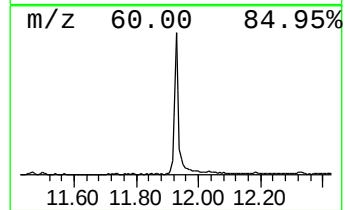
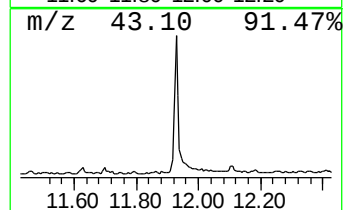
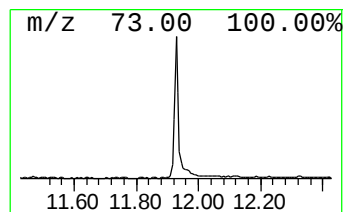
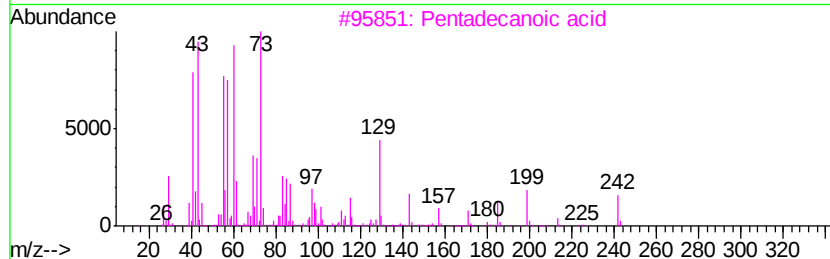
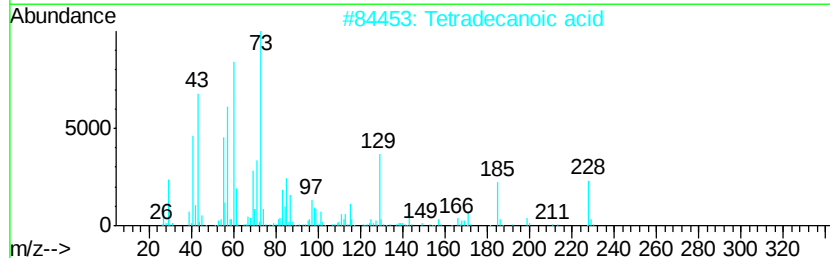
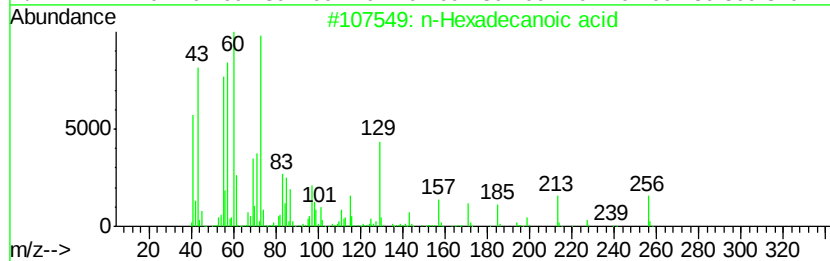
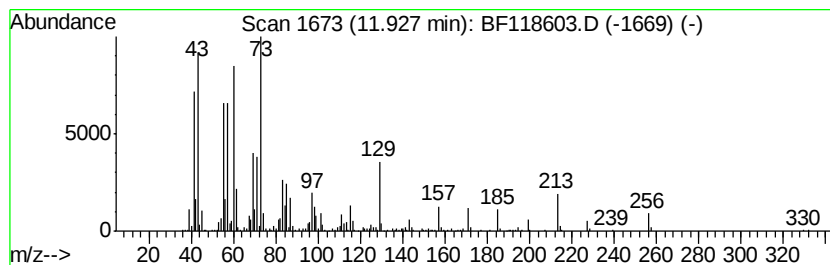
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	3.32 ng	743386	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4		Octadecanoic acid	284	C18H36O2	000057-11-4	76
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



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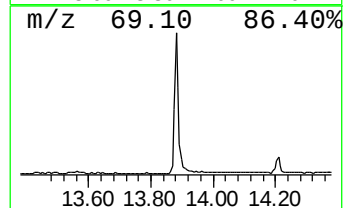
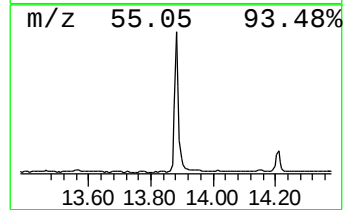
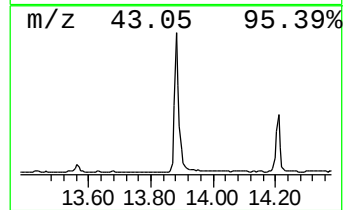
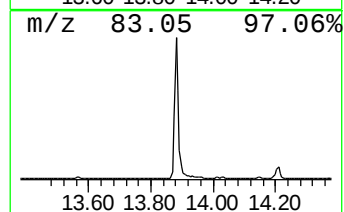
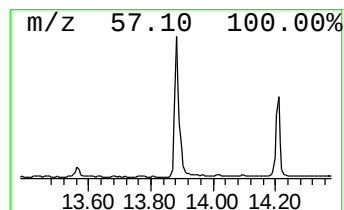
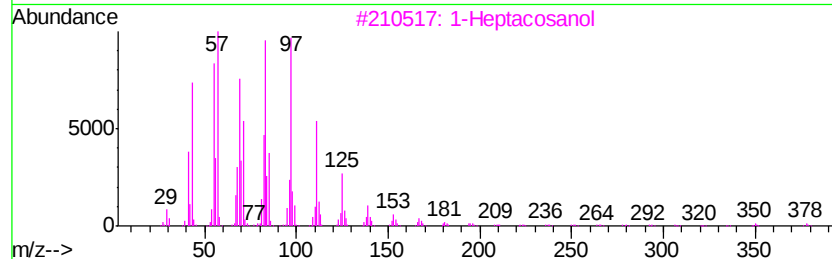
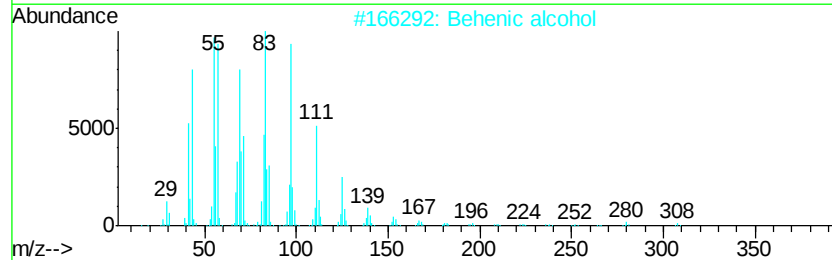
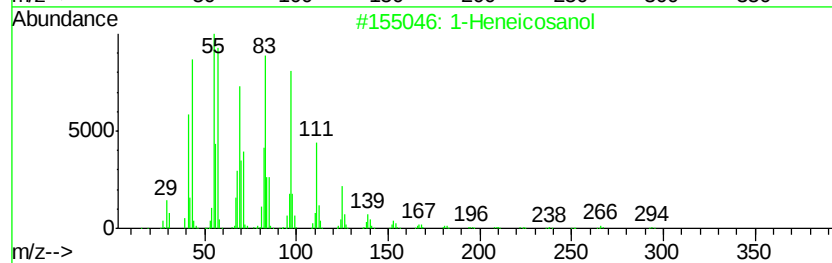
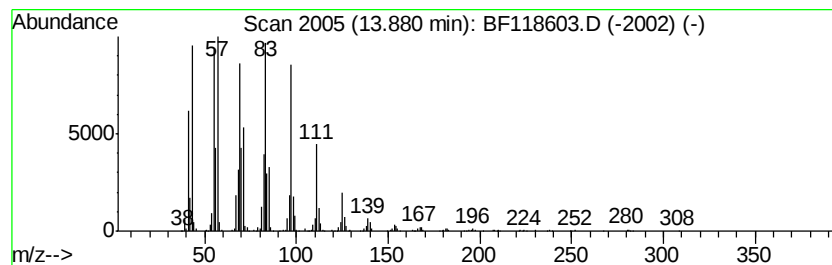
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Peak Number 6 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.88	12.29 ng	1844620	Chrysene-d12		14.03
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	1-Heptacosanol	396	C27H56O	002004-39-9	94
4	Cyclopentadecane	210	C15H30	000295-48-7	93
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93



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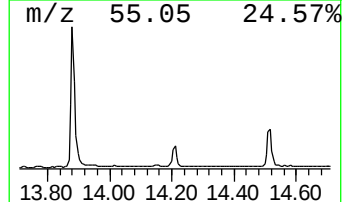
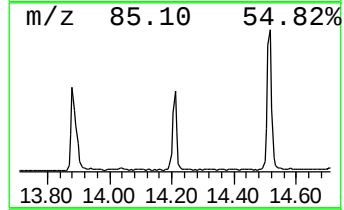
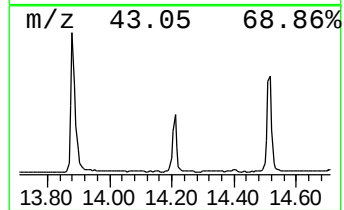
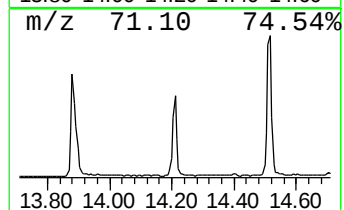
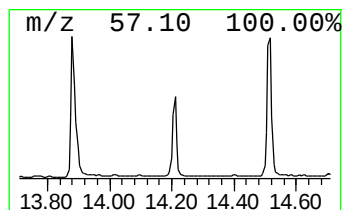
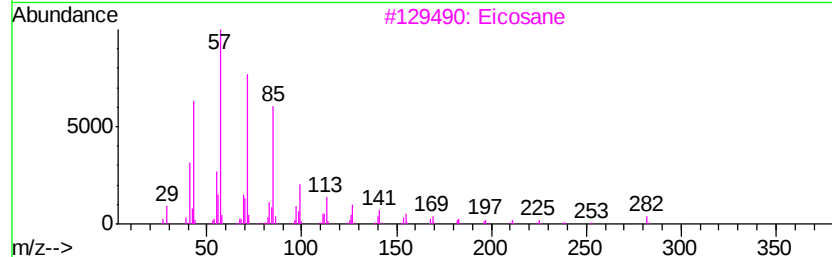
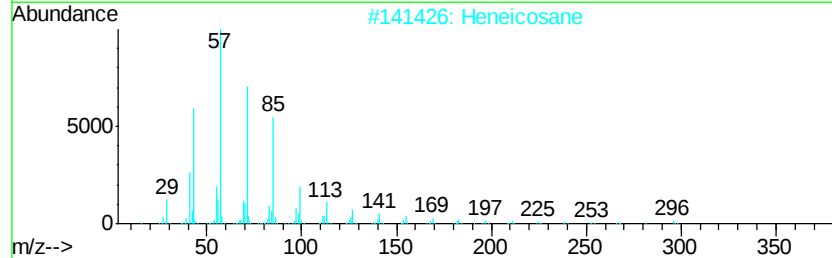
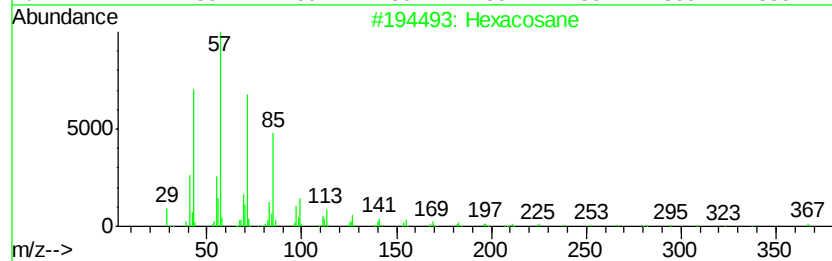
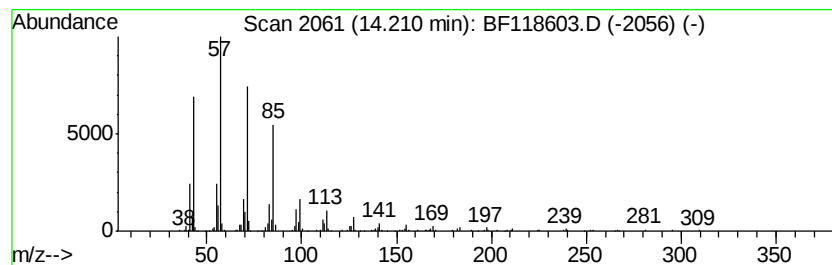
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	3.22 ng	483761	Chrysene-d12	14.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118603.D
Acq On : 20 Jan 2020 20:51
Operator : CG/JU
Sample : L1154-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3-0-15

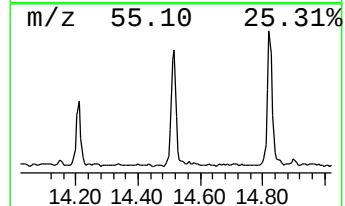
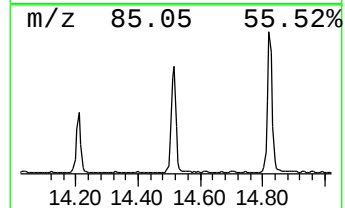
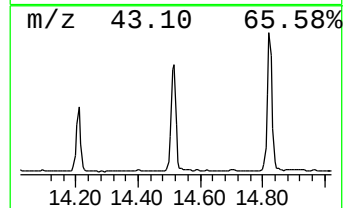
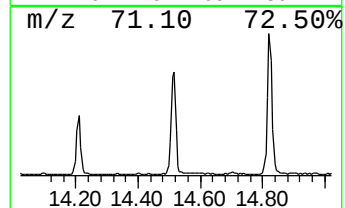
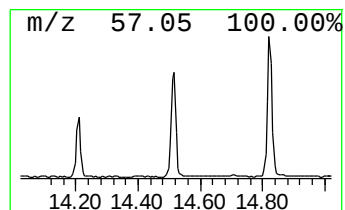
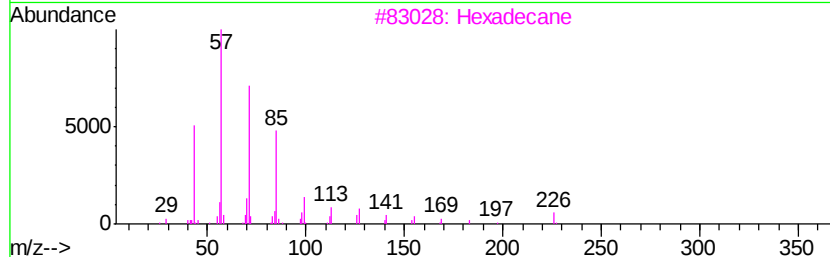
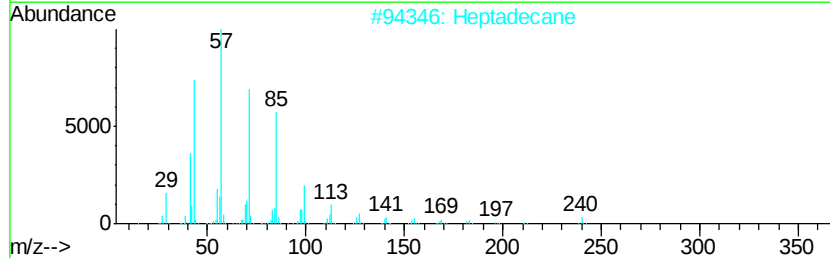
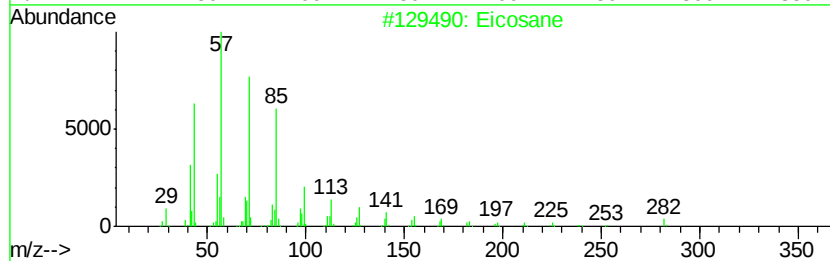
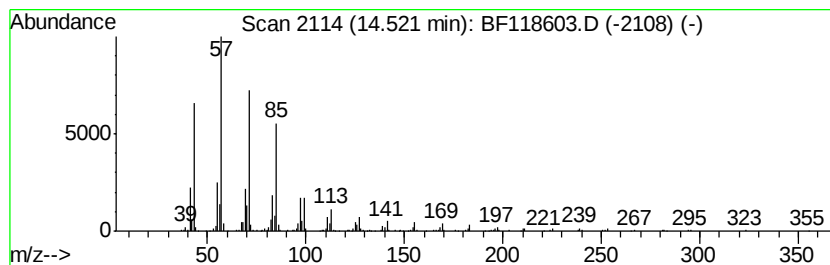
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	6.49 ng	973446	Chrysene-d12	14.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Heptadecane		240	C17H36	000629-78-7	93
3	Hexadecane		226	C16H34	000544-76-3	93
4	Hexacosane		366	C26H54	000630-01-3	91
5	Heneicosane		296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
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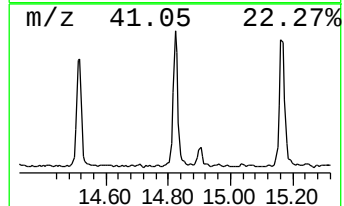
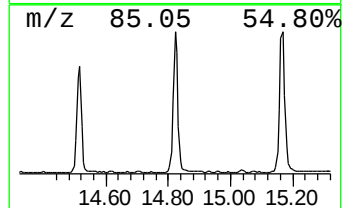
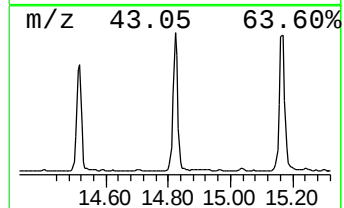
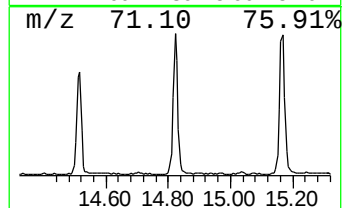
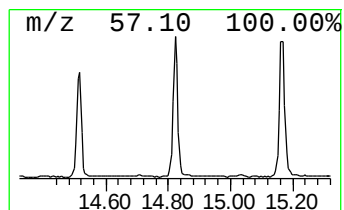
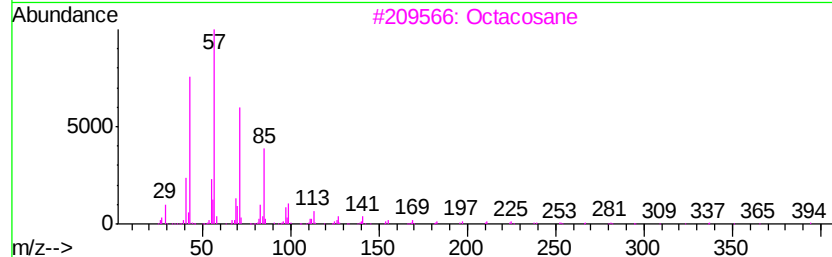
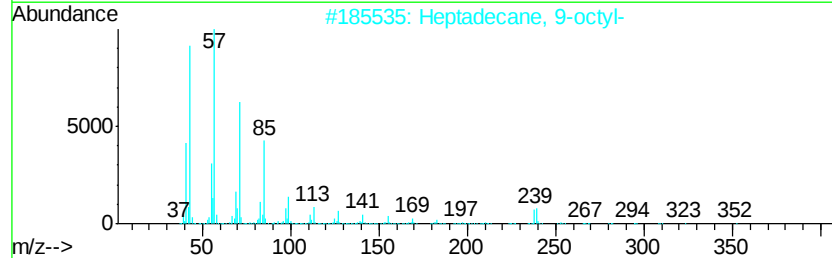
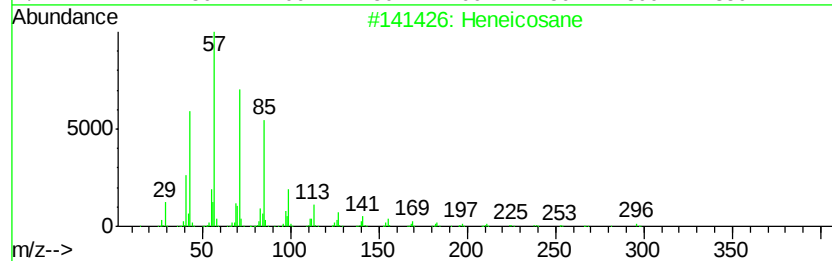
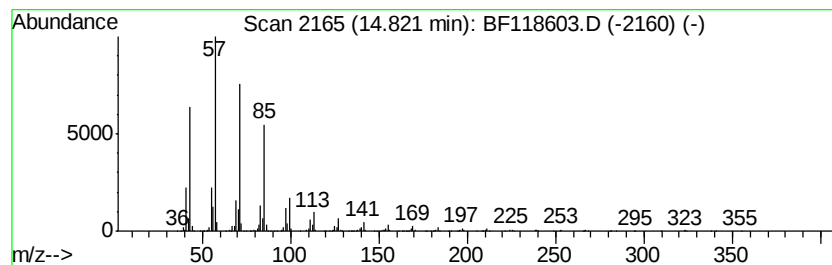
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heptadecane, 9-octyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	9.29 ng	1275830	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118603.D
Acq On : 20 Jan 2020 20:51
Operator : CG/JU
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Misc :
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ClientSampleId :
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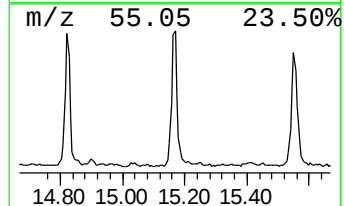
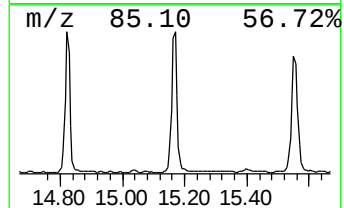
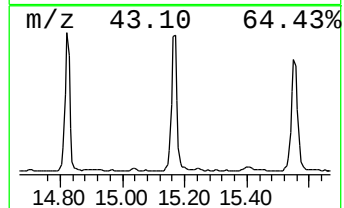
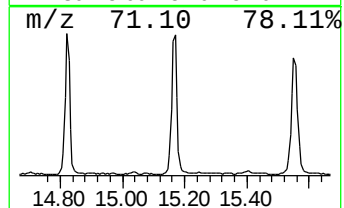
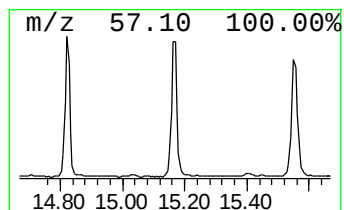
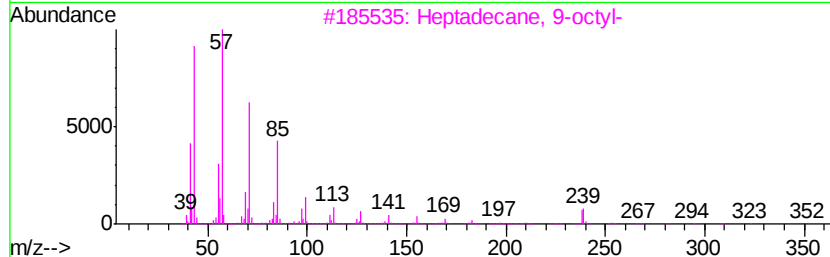
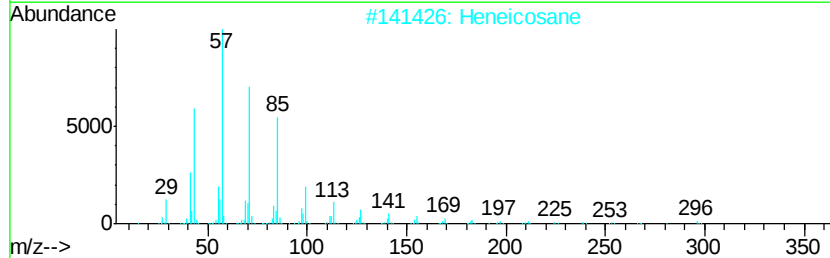
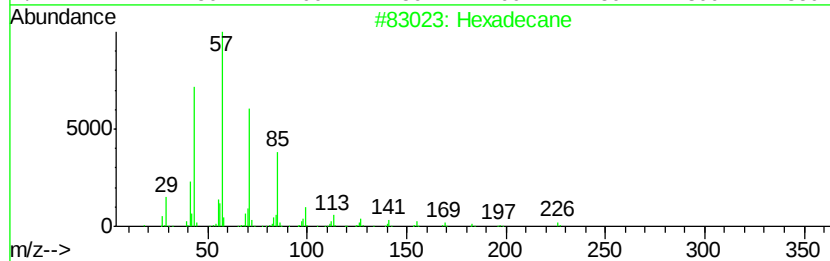
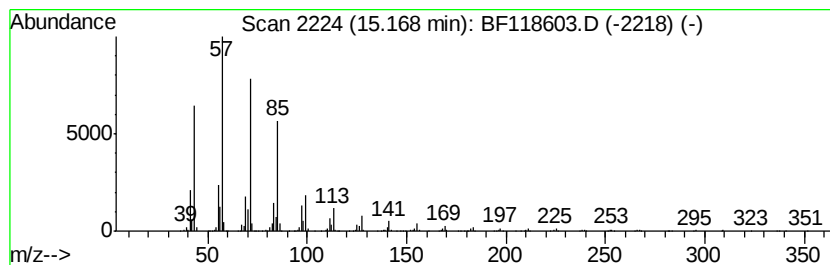
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	10.73 ng	1473660	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118603.D
Acq On : 20 Jan 2020 20:51
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Misc :
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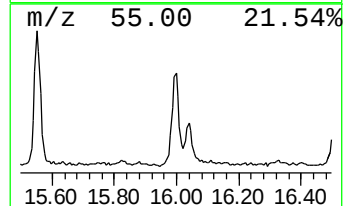
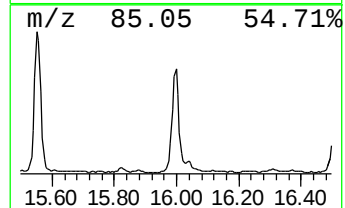
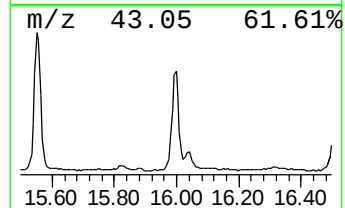
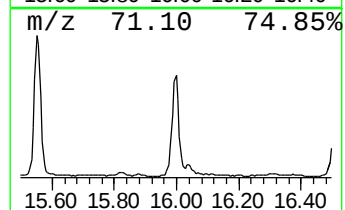
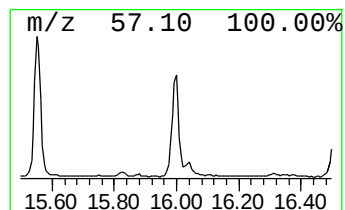
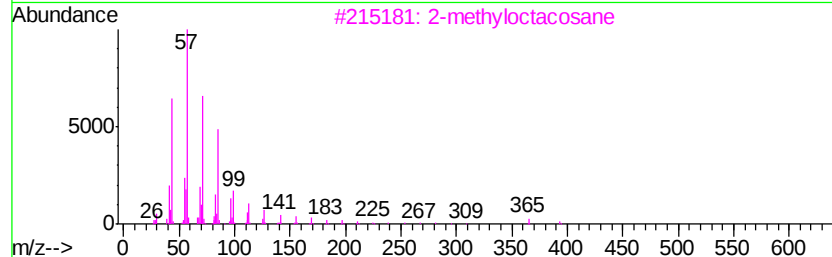
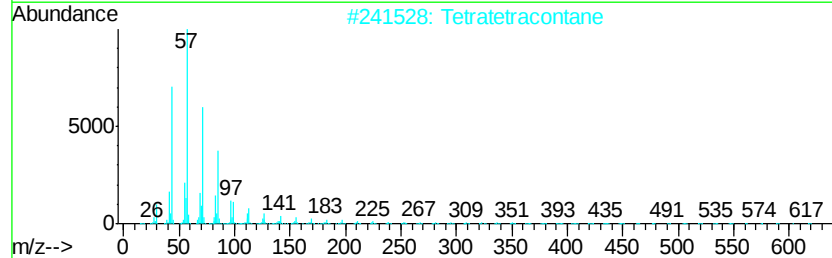
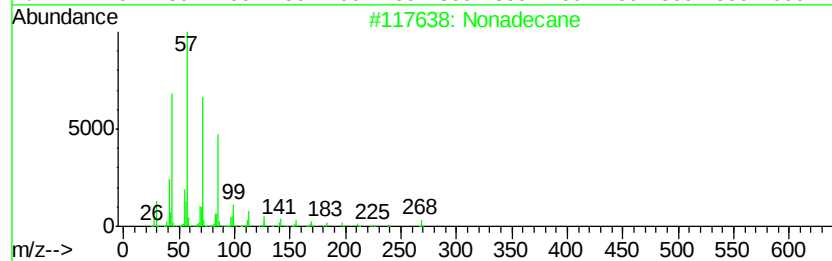
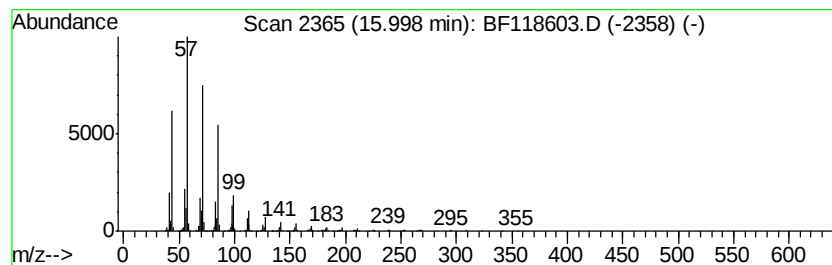
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	8.04 ng	1104190	Perylene-d12	15.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Tetratetracontane		619	C44H90	007098-22-8	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118603.D
Acq On : 20 Jan 2020 20:51
Operator : CG/JU
Sample : L1154-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3-0-15

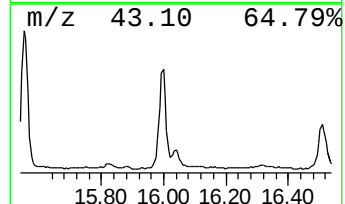
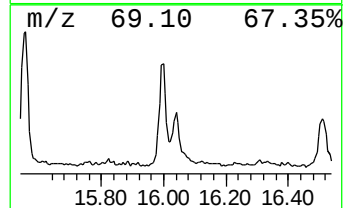
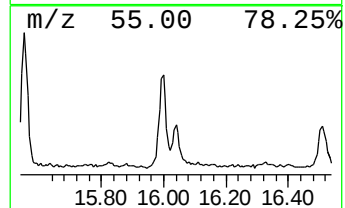
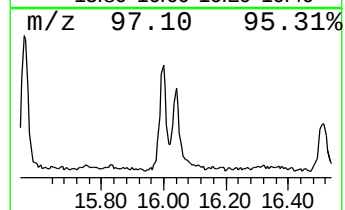
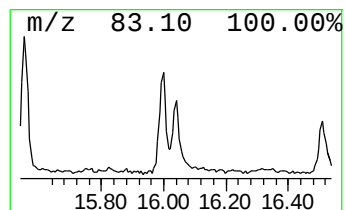
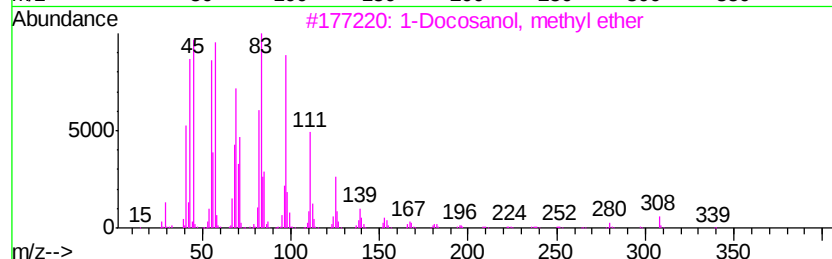
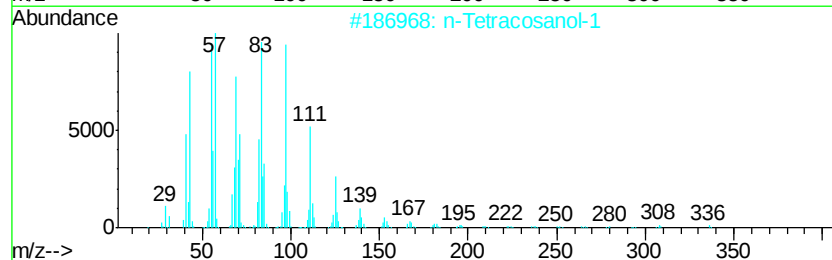
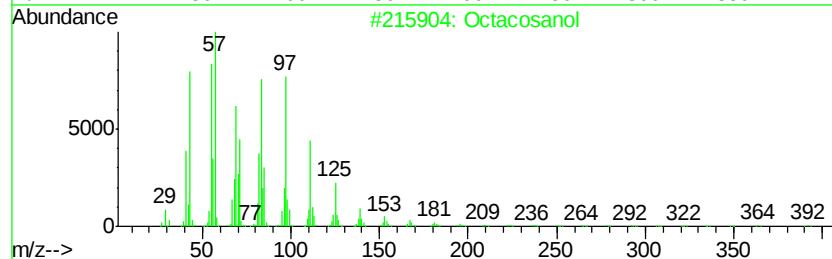
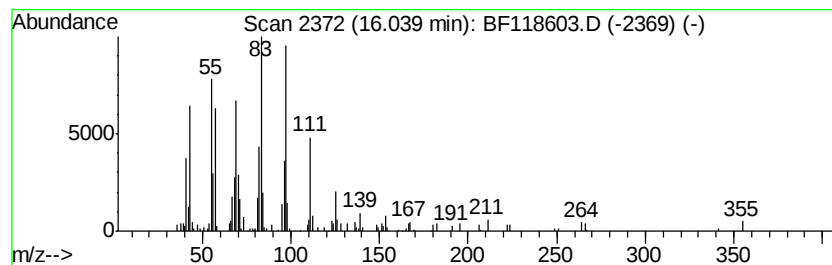
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.17 ng	297484	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosanol	410	C28H58O	000557-61-9	58
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	53
3			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	49
4			1-Heneicosanol	312	C21H44O	015594-90-8	46
5			9-Nonadecene	266	C19H38	031035-07-1	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012020\
Data File : BF118603.D
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ALS Vial : 13 Sample Multiplier: 1

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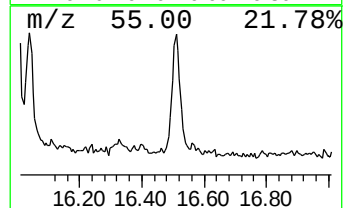
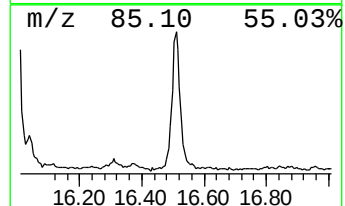
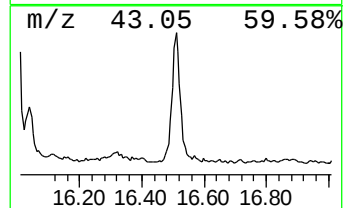
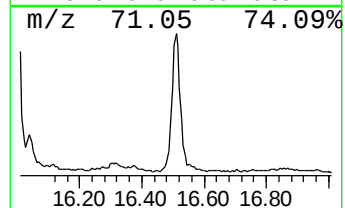
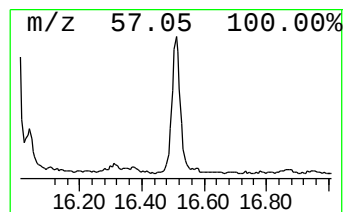
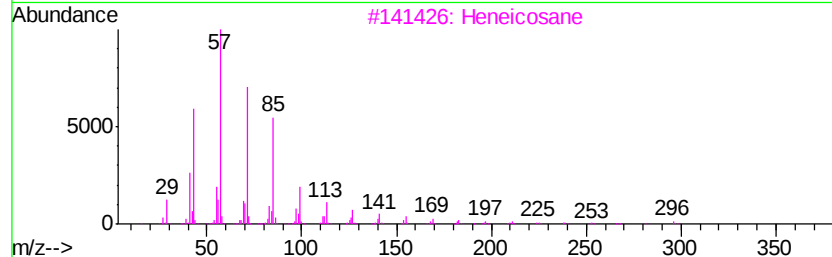
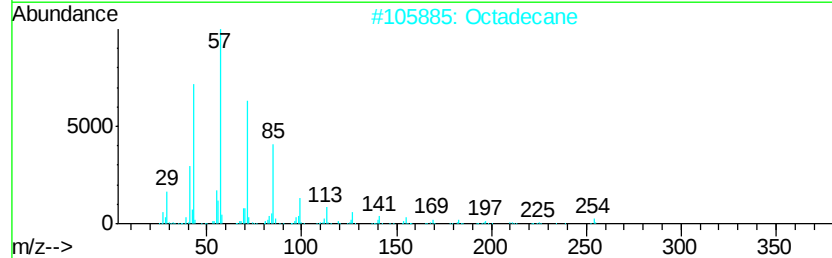
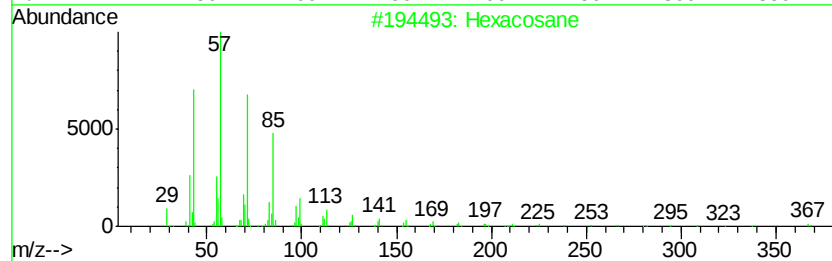
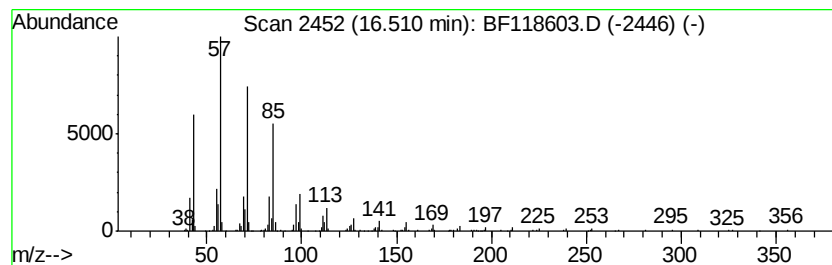
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 2-methyloctacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	4.13 ng	567658	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Octadecane	254	C18H38	000593-45-3	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane	394	C28H58	000630-02-4	91
5			2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012020\
 Data File : BF118603.D
 Acq On : 20 Jan 2020 20:51
 Operator : CG/JU
 Sample : L1154-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-0-15

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
Butane, 2-methoxy...	2.22	25.6	ng	3700420	1	6.88	2894390	20.0
2-Pentanone, 4-hy...	5.10	15.6	ng	2260400	1	6.88	2894390	20.0
unknown6.65	6.65	77.6	ng	11230100	1	6.88	2894390	20.0
n-Hexadecanoic acid	11.93	3.3	ng	743386	4	11.40	4478000	20.0
1-Heneicosanol	13.88	12.3	ng	1844620	5	14.03	3001530	20.0
Hexacosane	14.21	3.2	ng	483761	5	14.03	3001530	20.0
Eicosane	14.52	6.5	ng	973446	5	14.03	3001530	20.0
Heptadecane, 9-oc...	14.82	9.3	ng	1275830	6	15.50	2747830	20.0
Hexadecane	15.17	10.7	ng	1473660	6	15.50	2747830	20.0
Nonadecane	16.00	8.0	ng	1104190	6	15.50	2747830	20.0
Octacosanol	16.04	2.2	ng	297484	6	15.50	2747830	20.0
2-methyloctacosane	16.51	4.1	ng	567658	6	15.50	2747830	20.0