

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
 Data File : BF119703.D
 Acq On : 2 Apr 2020 8:01
 Operator : MA/SJ
 Sample : L1752-05
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-02-033020

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-BF031820.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	2.210	17	21	27	rVB	119170	158579	2.66%	0.324%
2	5.034	497	501	508	rBV	199224	261335	4.39%	0.533%
3	5.434	563	569	572	rBV	5917743	5957004	100.00%	12.158%
4	6.445	735	741	744	rBV	5398699	5841874	98.07%	11.923%
5	6.639	771	774	779	rBV	85264	80812	1.36%	0.165%
6	6.816	800	804	807	rBV	1826304	1572838	26.40%	3.210%
7	6.969	826	830	834	rBV	4768603	4662852	78.28%	9.517%
8	7.381	894	900	903	rBV	3745796	3841188	64.48%	7.840%
9	8.098	1016	1022	1025	rBV	2503095	2005297	33.66%	4.093%
10	9.180	1200	1206	1209	rBV	6324449	5936519	99.66%	12.116%
11	9.257	1216	1219	1223	rBV2	69135	80432	1.35%	0.164%
12	9.569	1268	1272	1275	rBV	1055582	793429	13.32%	1.619%
13	9.786	1307	1309	1314	rVB	73181	65617	1.10%	0.134%
14	9.857	1317	1321	1324	rVB	2438206	2023956	33.98%	4.131%
15	10.645	1450	1455	1458	rBV	3789433	3311650	55.59%	6.759%
16	10.763	1472	1475	1480	rBV	79180	84066	1.41%	0.172%
17	11.210	1548	1551	1554	rBV2	69303	63358	1.06%	0.129%
18	11.339	1569	1573	1576	rBV	1905953	1762804	29.59%	3.598%
19	11.363	1576	1577	1580	rVB	117072	70637	1.19%	0.144%
20	11.739	1638	1641	1644	rBV	183918	143851	2.41%	0.294%
21	11.869	1660	1663	1670	rBV	480987	452127	7.59%	0.923%
22	12.421	1754	1757	1759	rBV2	424625	416118	6.99%	0.849%
23	12.445	1759	1761	1764	rVB	775503	579123	9.72%	1.182%
24	12.533	1773	1776	1777	rBV	90942	73329	1.23%	0.150%
25	12.551	1777	1779	1782	rVB	105501	99753	1.67%	0.204%
26	12.657	1793	1797	1804	rBV	207370	225559	3.79%	0.460%
27	12.780	1811	1818	1821	rBV	111241	140947	2.37%	0.288%
28	12.927	1839	1843	1847	rBV	4129054	3847563	64.59%	7.853%
29	13.163	1879	1883	1890	rBV2	64021	111232	1.87%	0.227%
30	13.410	1922	1925	1930	rBV3	67396	66266	1.11%	0.135%
31	13.833	1993	1997	2007	rBV	391797	531773	8.93%	1.085%
32	13.980	2018	2022	2025	rBV	1398497	1251929	21.02%	2.555%
33	14.151	2049	2051	2055	rVB	115856	98345	1.65%	0.201%
34	14.457	2100	2103	2113	rBV3	220356	402244	6.75%	0.821%

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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-BF031820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.762	2152	2155	2159	rBV	195664	207854	3.49%	0.424%
36	15.098	2209	2212	2215	rBV	232497	258285	4.34%	0.527%
37	15.439	2266	2270	2274	rBV	919065	1091099	18.32%	2.227%
38	15.480	2274	2277	2282	rVB	173272	217173	3.65%	0.443%
39	15.909	2347	2350	2356	rVB	143786	208727	3.50%	0.426%

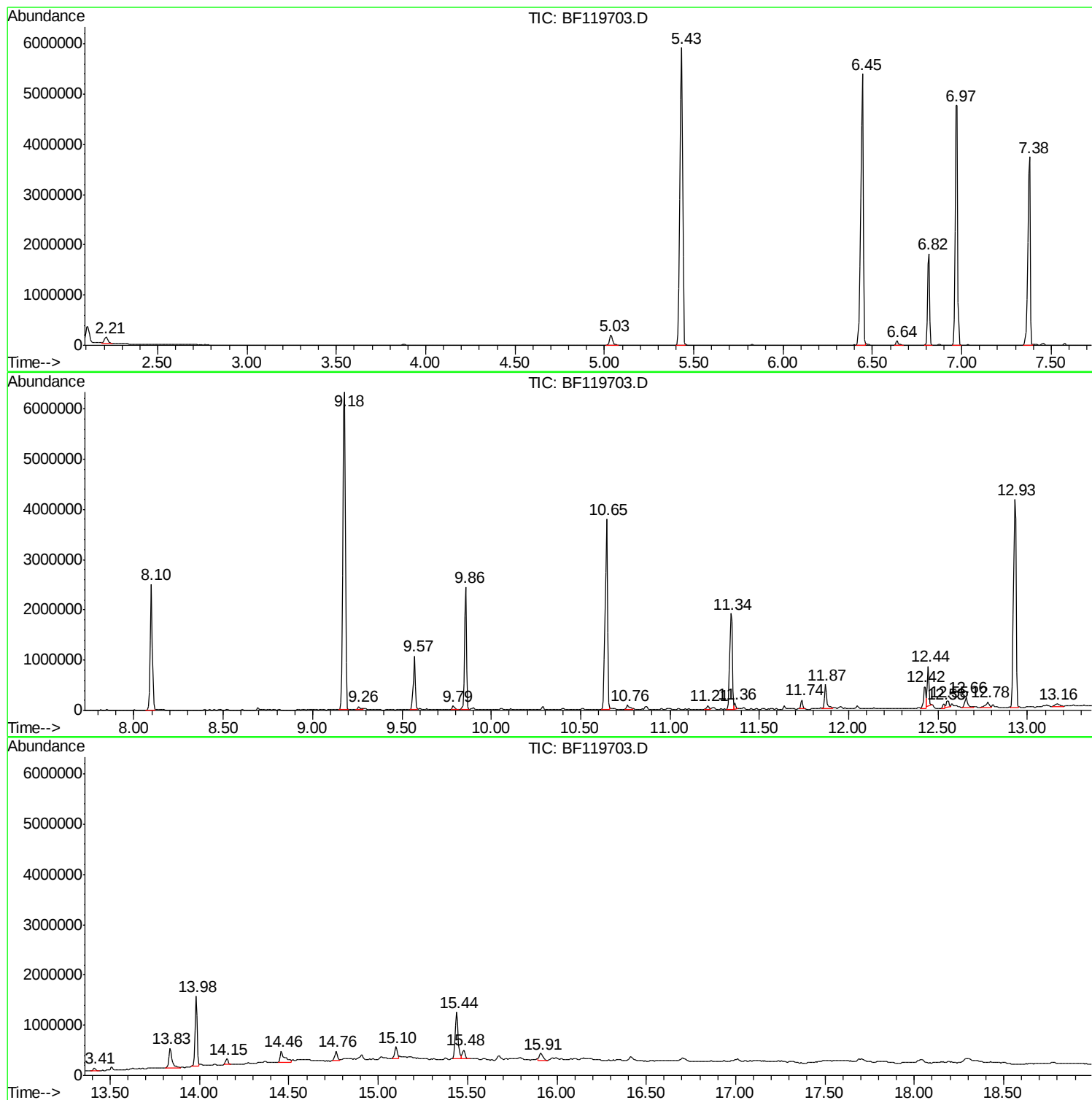
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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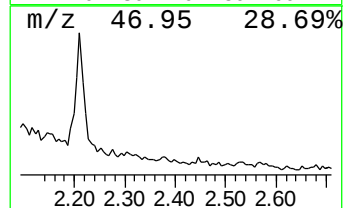
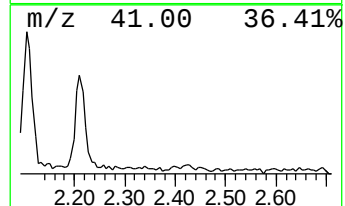
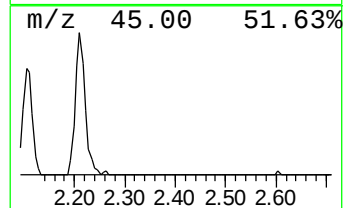
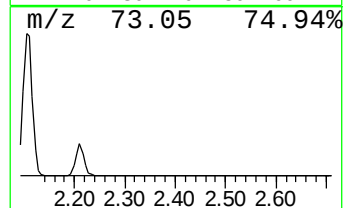
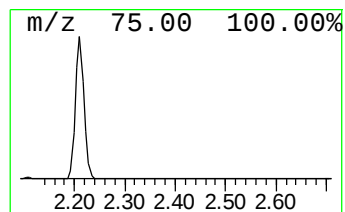
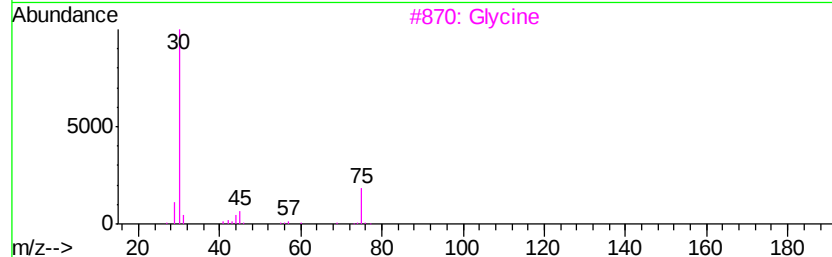
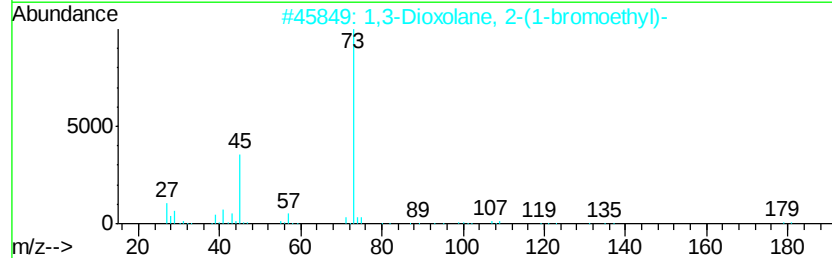
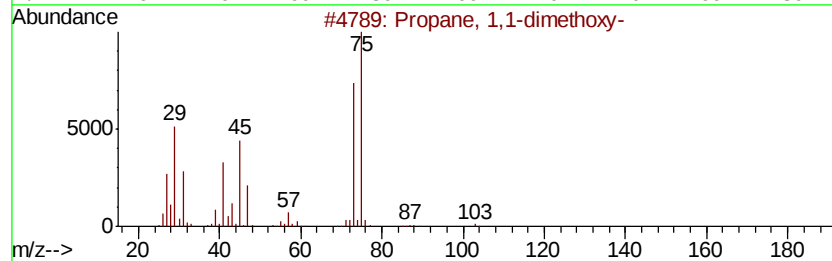
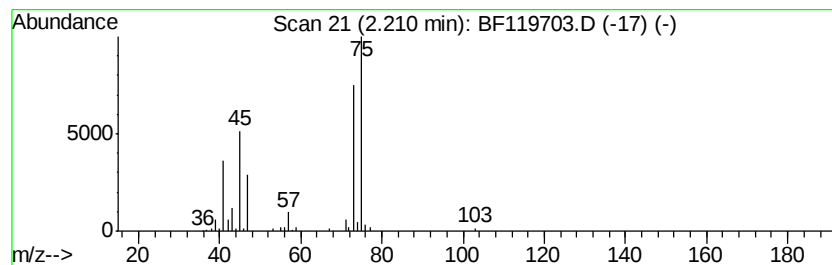
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.21	2.02 ng	158579	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2		1,3-Dioxolane, 2-(1-bromoethyl)-	180	C5H9BrO2	005267-73-2	16
3		Glycine	75	C2H5NO2	000056-40-6	9
4		Ethanol, 2-methoxy-	76	C3H8O2	000109-86-4	9
5		Thiocyanic acid, methyl ester	73	C2H3NS	000556-64-9	9



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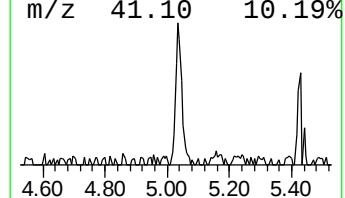
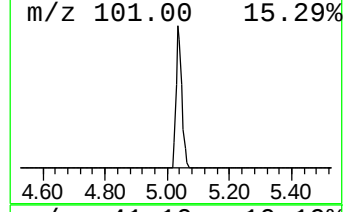
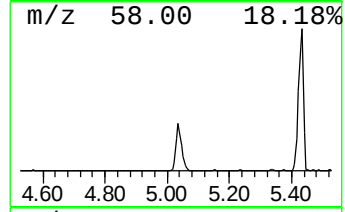
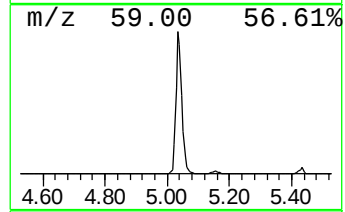
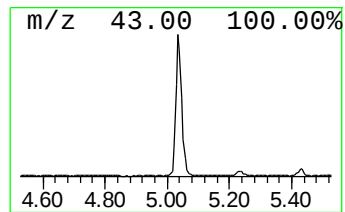
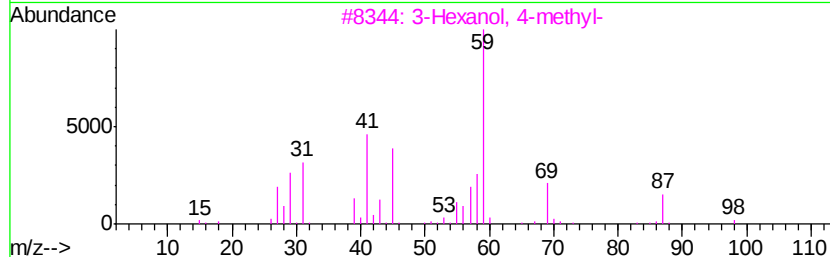
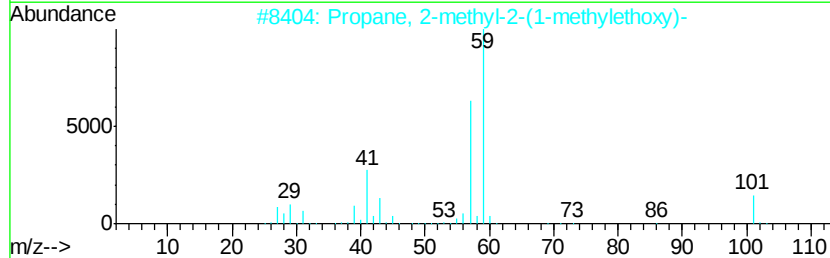
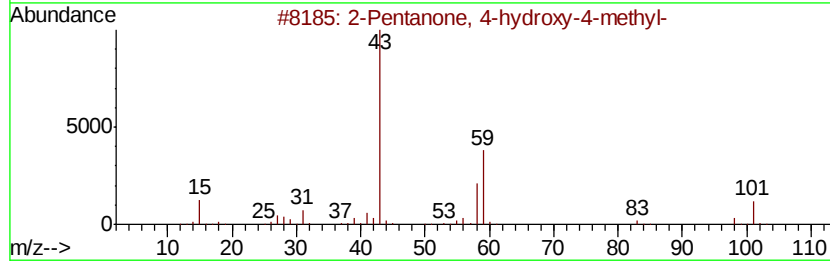
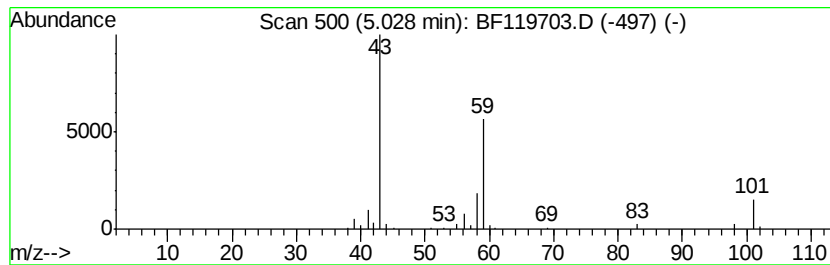
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	3.32 ng	261335	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		3-Octanol	130	C8H18O	000589-98-0	33
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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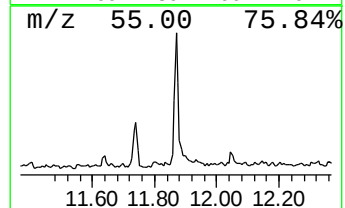
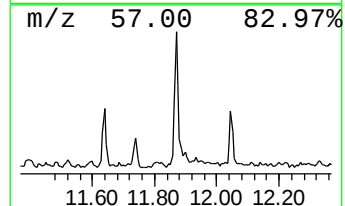
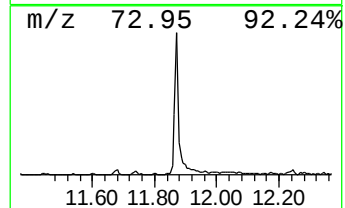
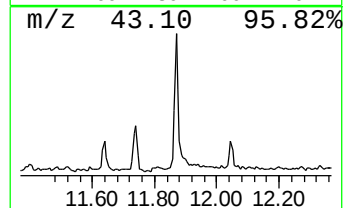
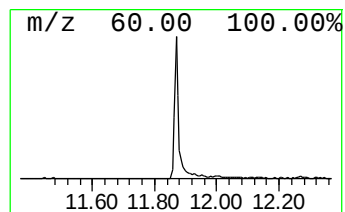
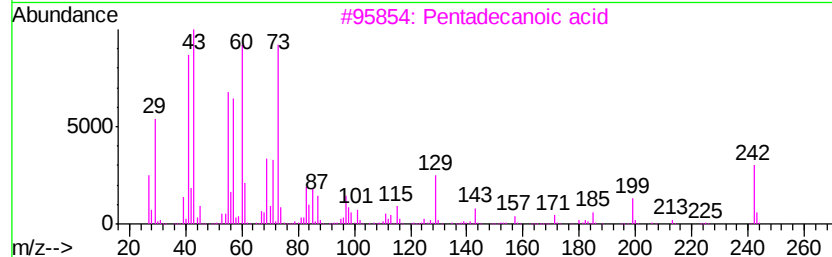
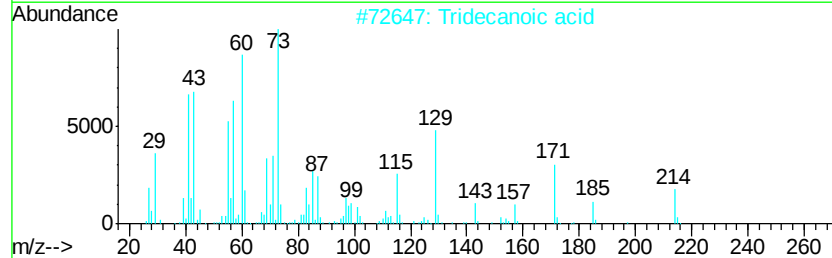
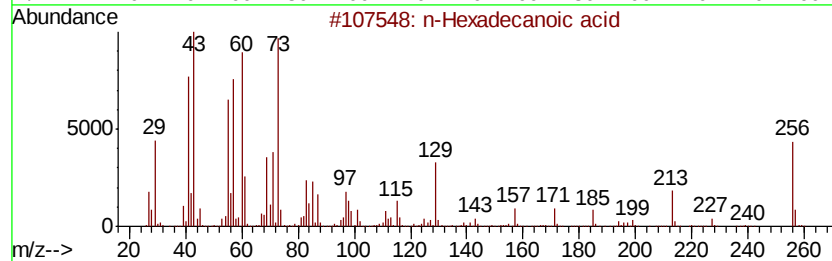
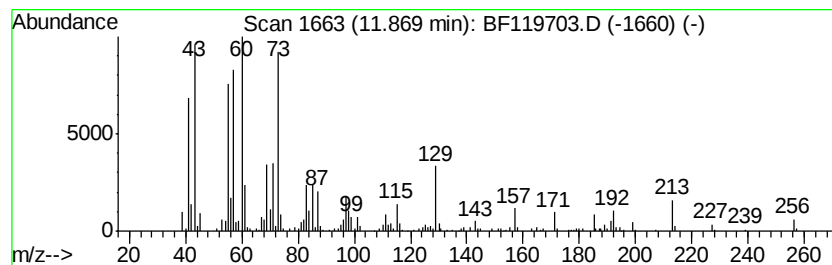
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	5.13 ng	452127	Phenanthrene-d10	11.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Undecanoic acid	186	C11H22O2	000112-37-8	59



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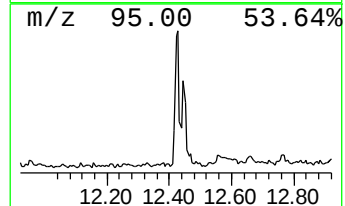
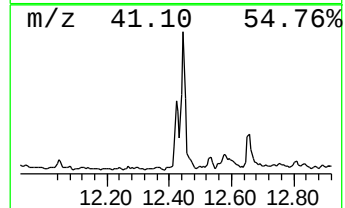
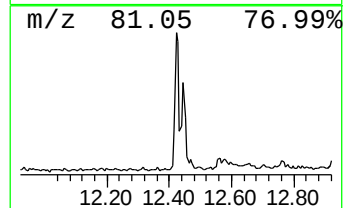
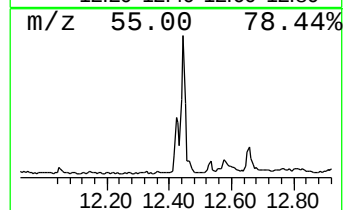
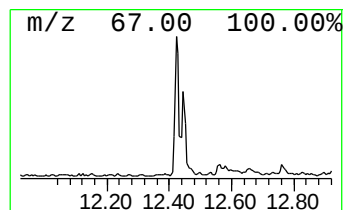
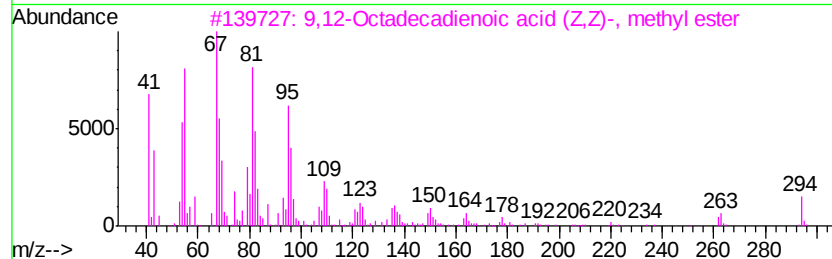
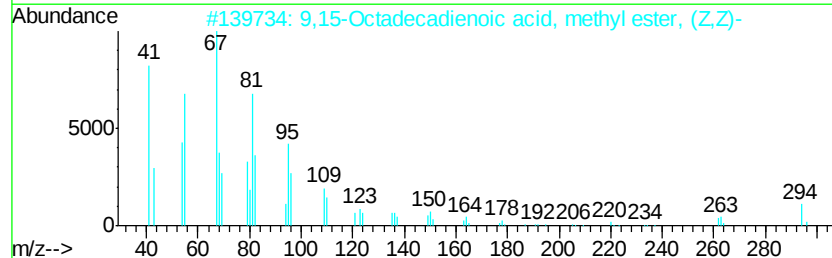
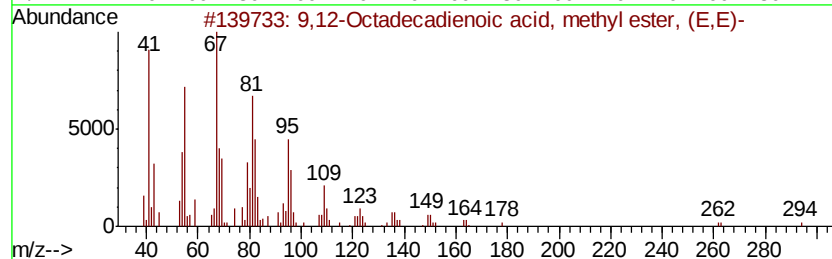
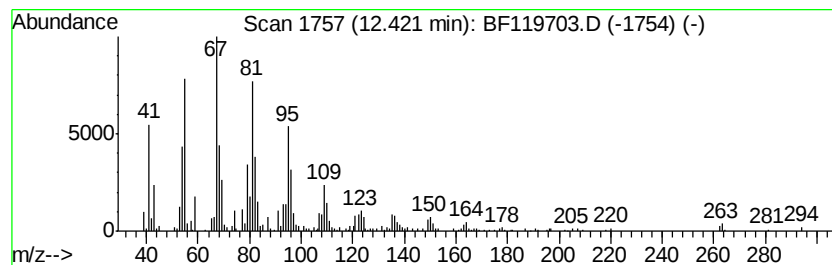
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9,12-Octadecadienoic acid, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.42	4.72 ng	416118	Phenanthrene-d10	11.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	98
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	97
5	11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	91



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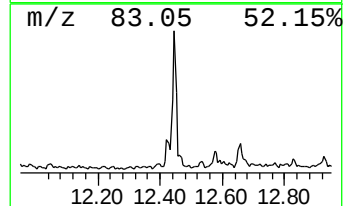
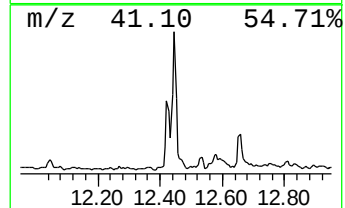
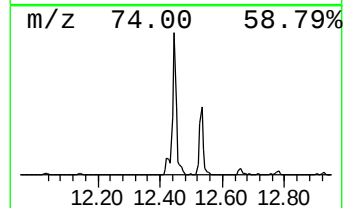
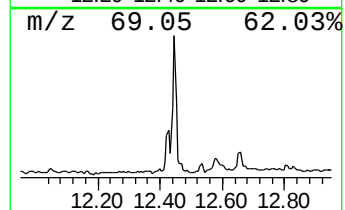
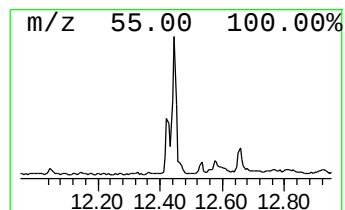
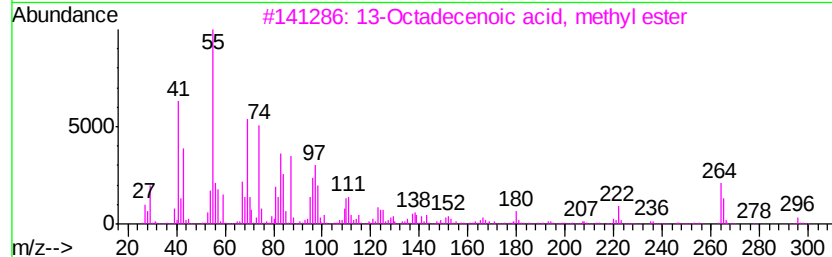
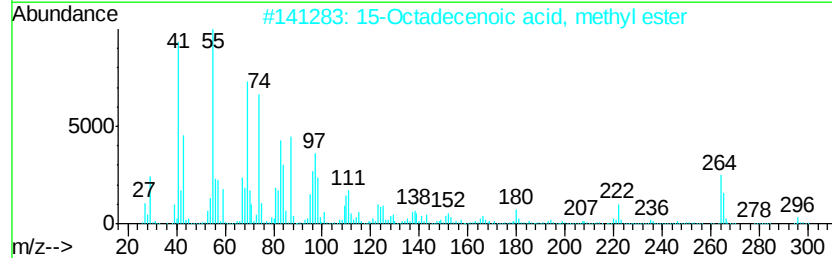
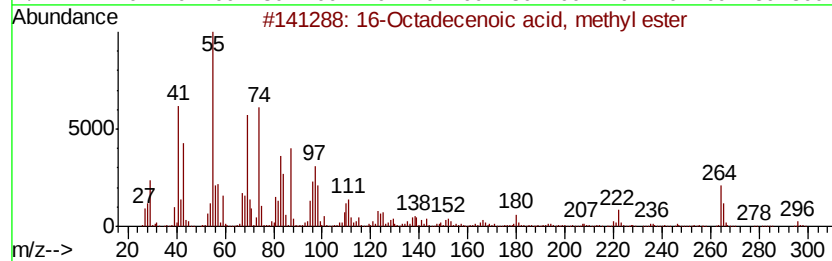
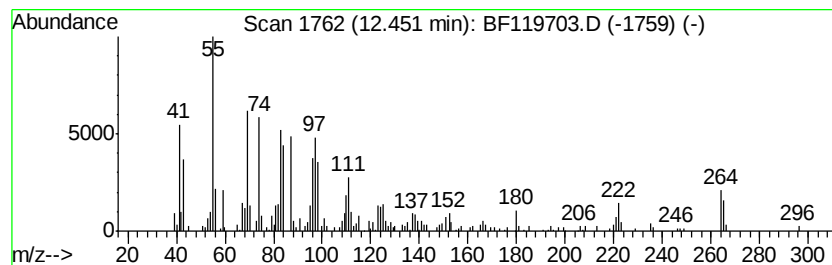
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 16-Octadecenoic acid, methyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.45	6.57 ng	579123	Phenanthrene-d10	11.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	98
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	98
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	96
4	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	96
5	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
Data File : BF119703.D
Acq On : 2 Apr 2020 8:01
Operator : MA/SJ
Sample : L1752-05
Misc :
ALS Vial : 53 Sample Multiplier: 1

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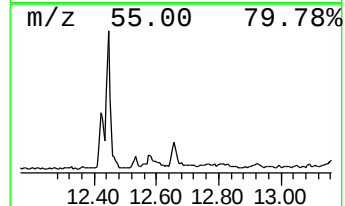
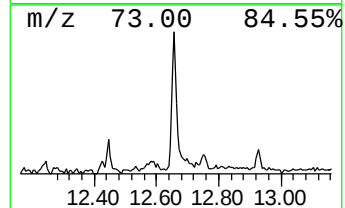
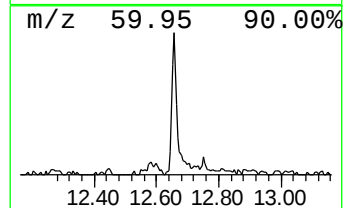
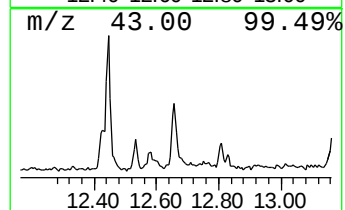
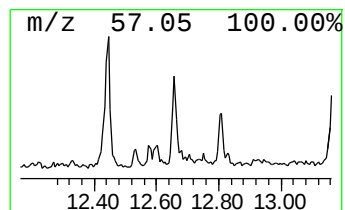
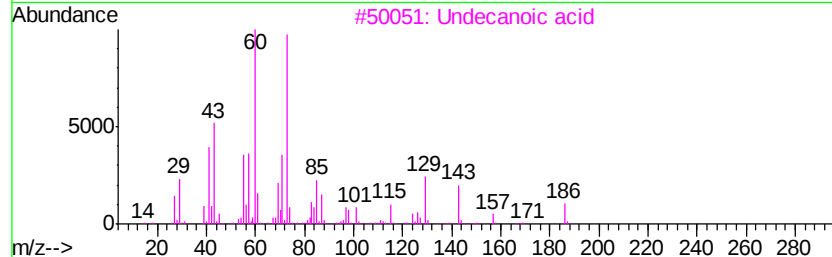
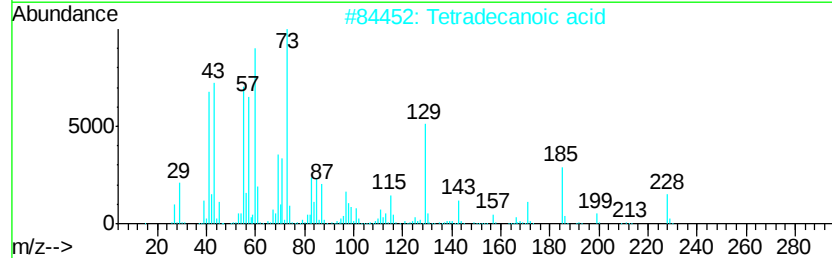
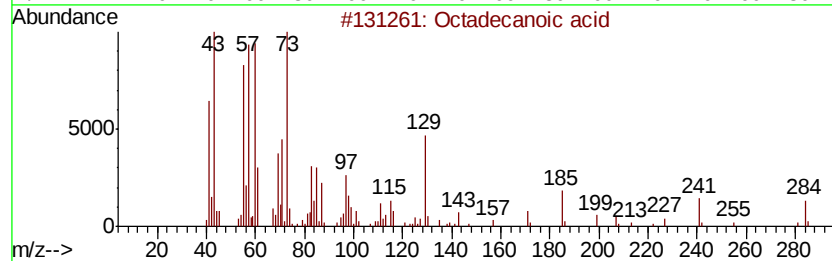
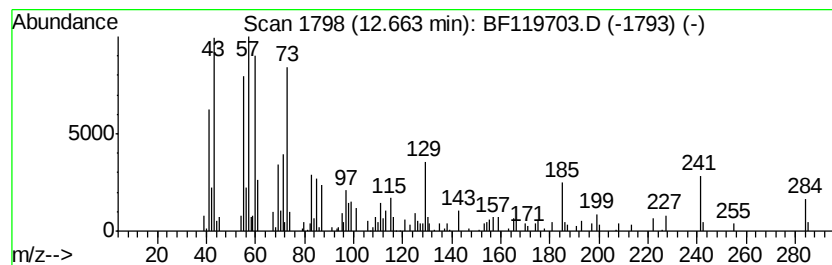
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	2.56 ng	225559	Phenanthrene-d10	11.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		Undecanoic acid	186	C11H22O2	000112-37-8	76
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
 Data File : BF119703.D
 Acq On : 2 Apr 2020 8:01
 Operator : MA/SJ
 Sample : L1752-05
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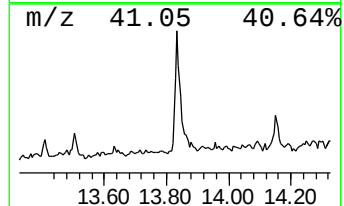
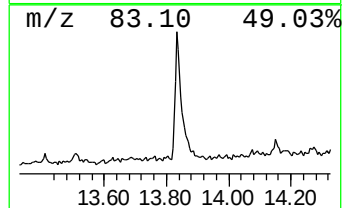
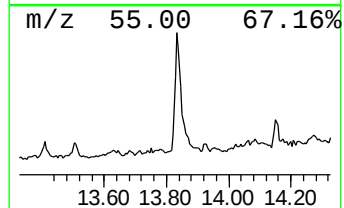
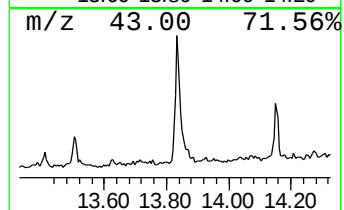
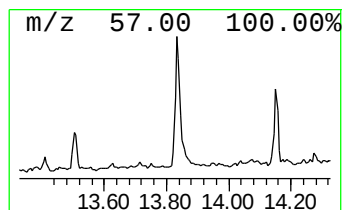
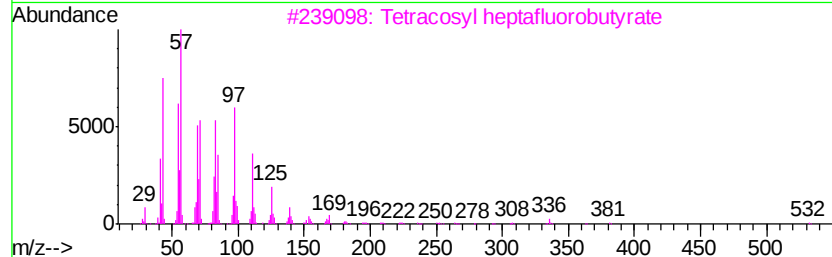
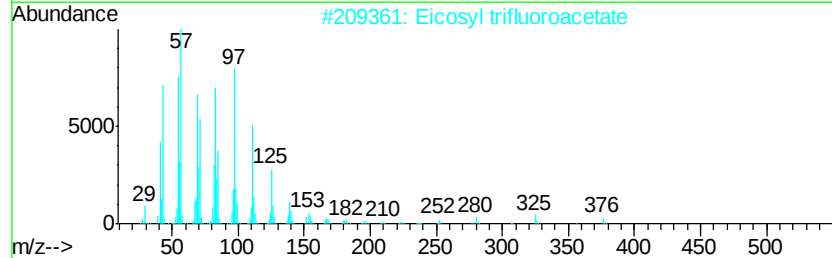
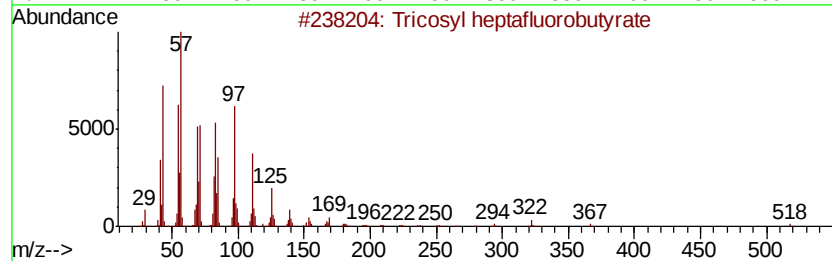
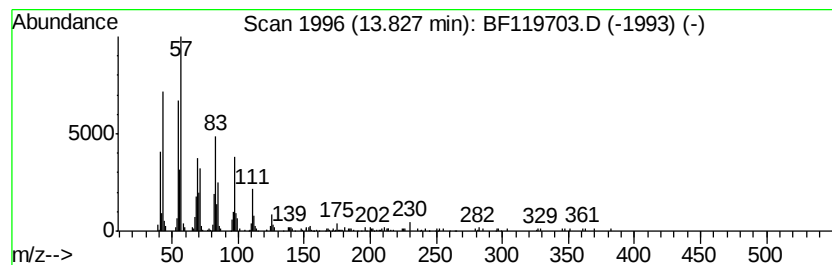
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Tricosyl heptafluorobutyrate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	8.50 ng	531773	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	91
2		Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91
3		Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
5		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
Data File : BF119703.D
Acq On : 2 Apr 2020 8:01
Operator : MA/SJ
Sample : L1752-05
Misc :
ALS Vial : 53 Sample Multiplier: 1

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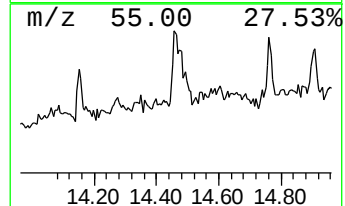
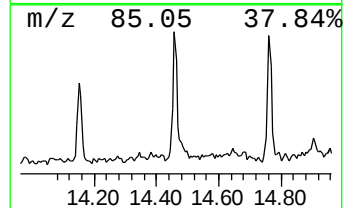
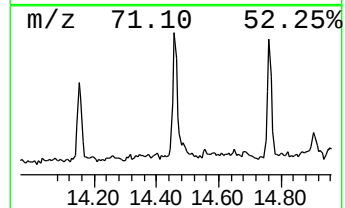
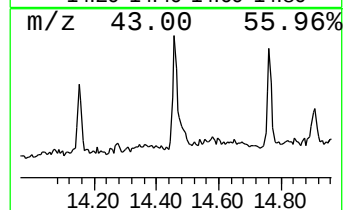
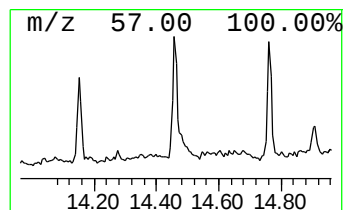
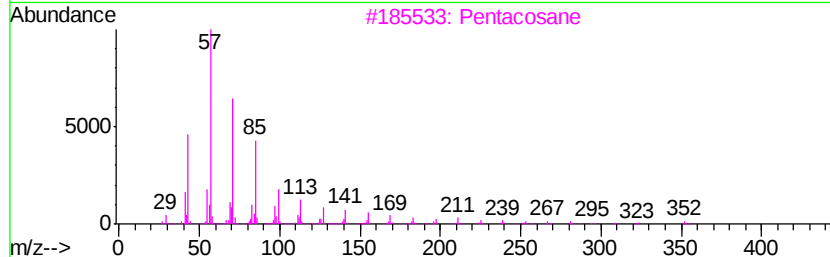
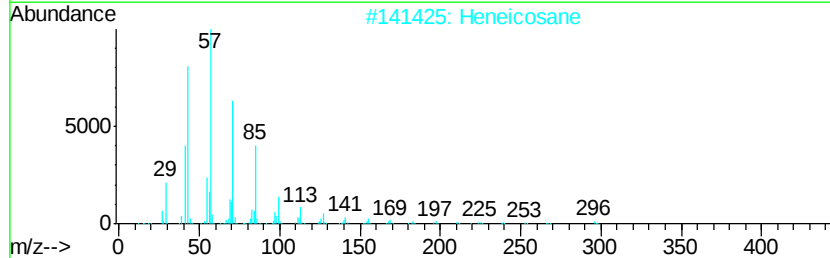
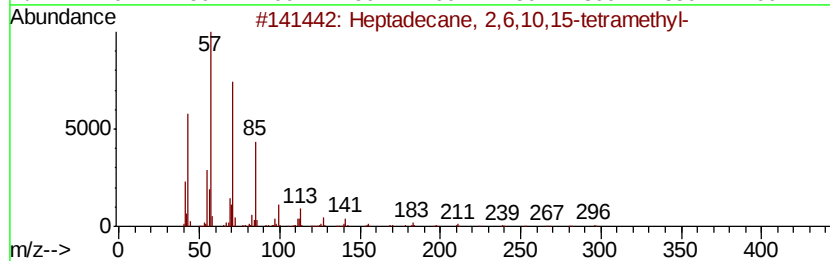
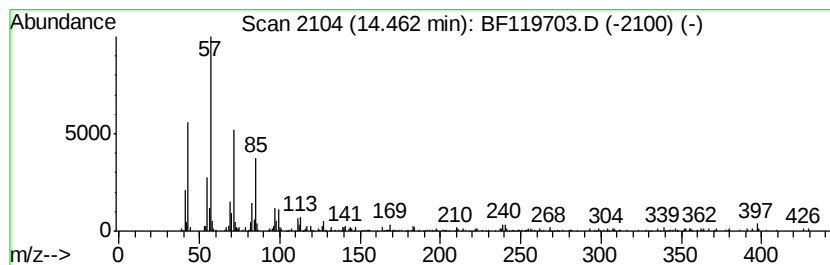
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 2,6,10,15-tetr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	6.43 ng	402244	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	98
2		Heneicosane	296	C21H44	000629-94-7	93
3		Pentacosane	352	C25H52	000629-99-2	93
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
Data File : BF119703.D
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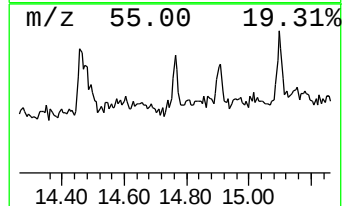
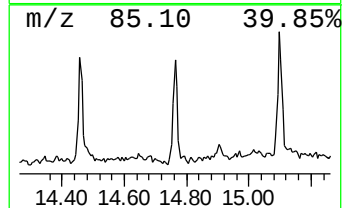
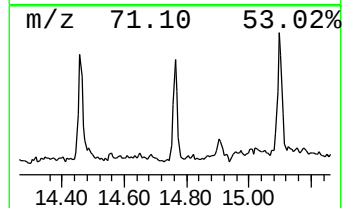
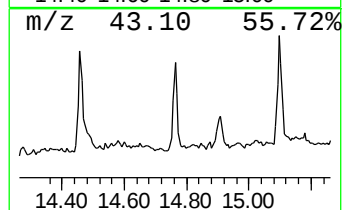
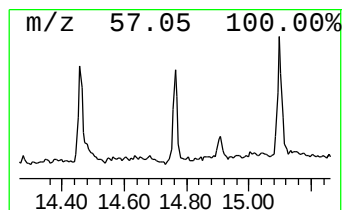
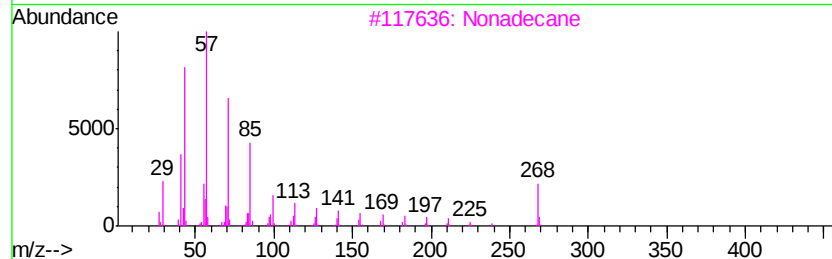
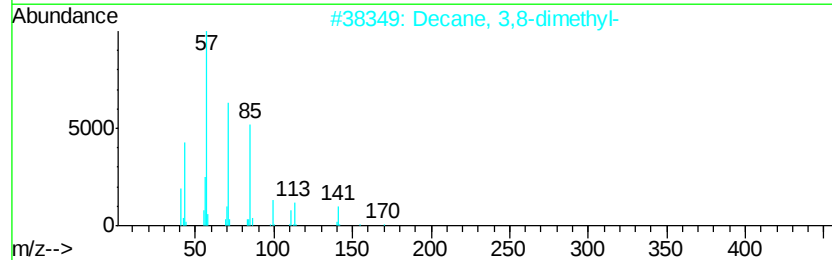
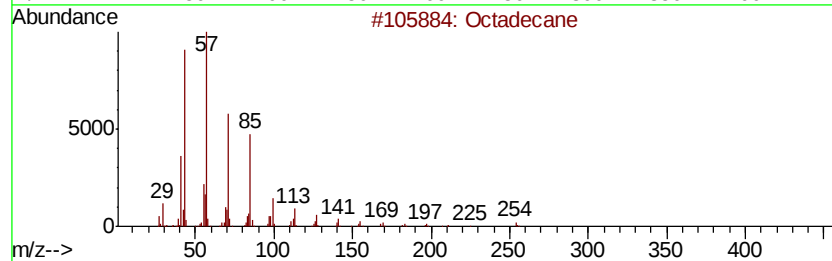
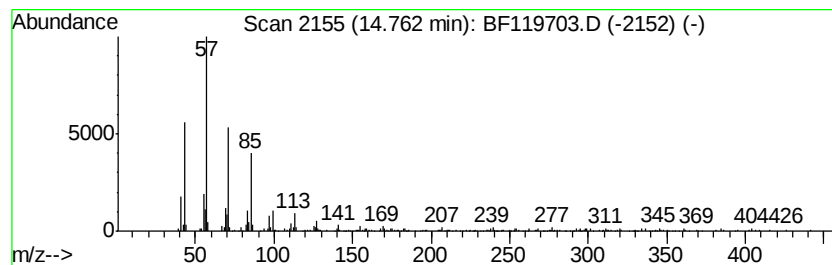
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	3.81 ng	207854	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	90
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87
3			Nonadecane	268	C19H40	000629-92-5	87
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
5			Tridecane, 7-propyl-	226	C16H34	055045-09-5	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
Data File : BF119703.D
Acq On : 2 Apr 2020 8:01
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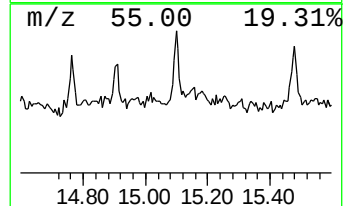
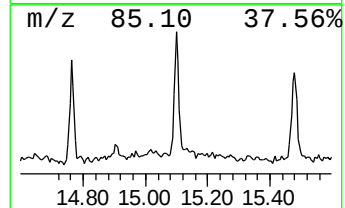
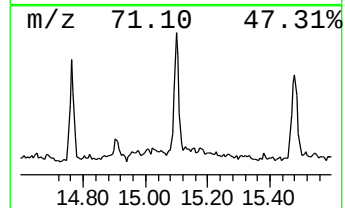
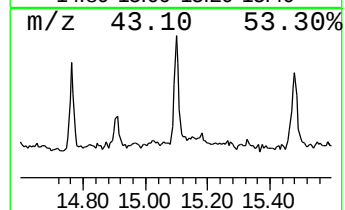
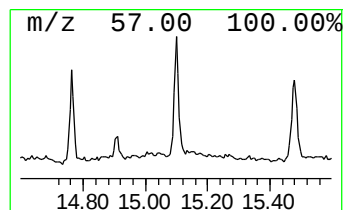
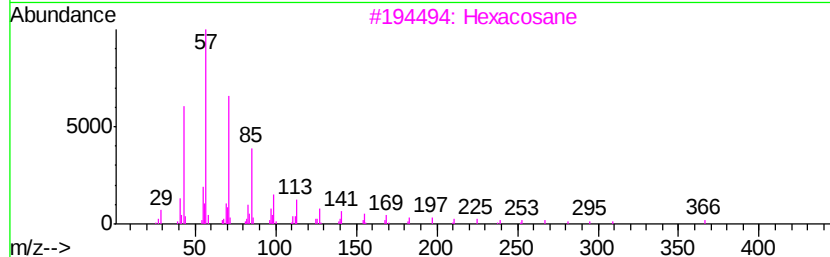
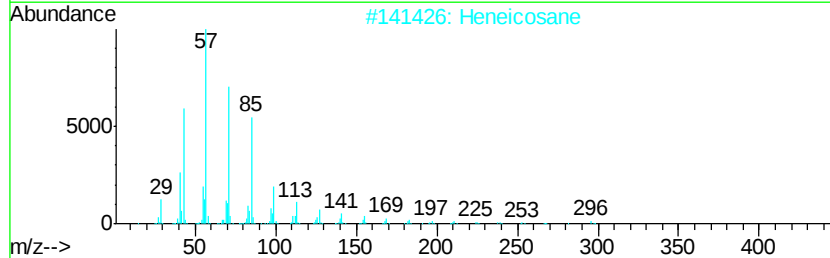
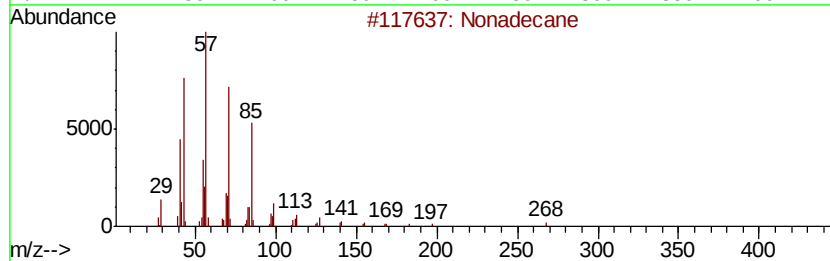
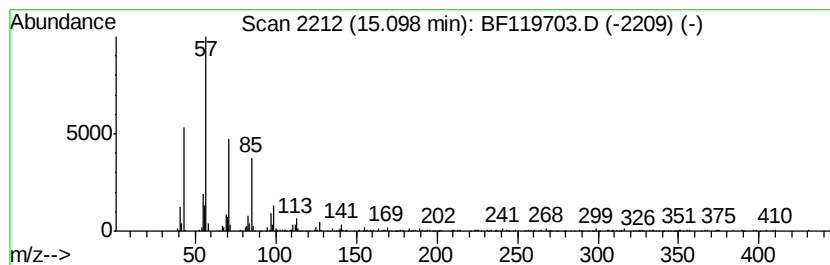
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	4.73 ng	258285	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	90
2		Heneicosane	296	C21H44	000629-94-7	81
3		Hexacosane	366	C26H54	000630-01-3	80
4		2-methyloctacosane	408	C29H60	1000376-72-8	74
5		Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
Data File : BF119703.D
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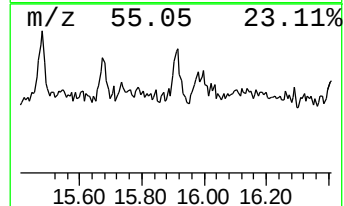
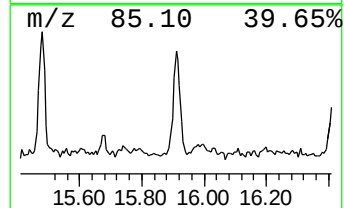
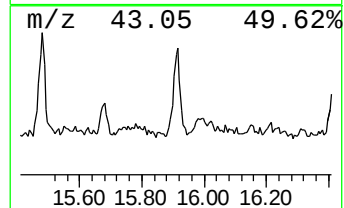
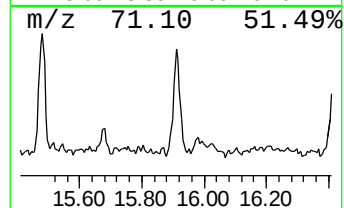
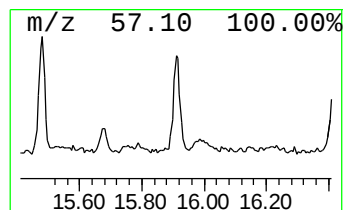
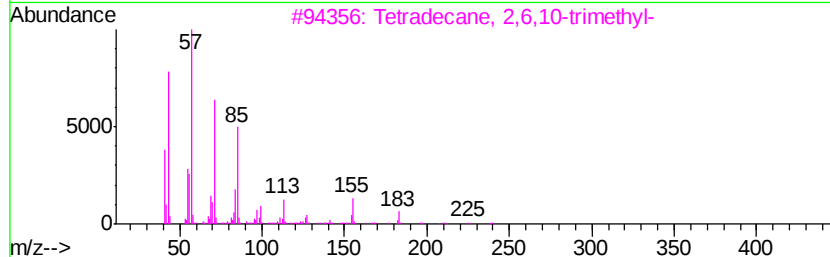
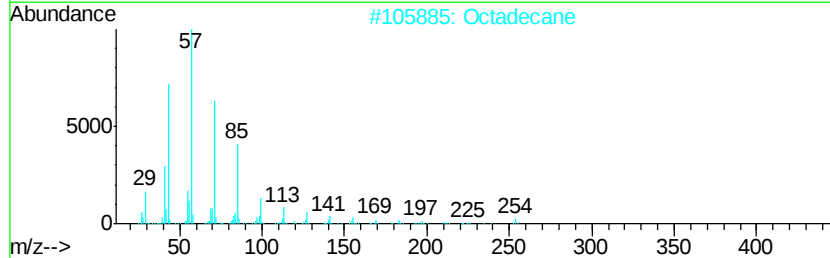
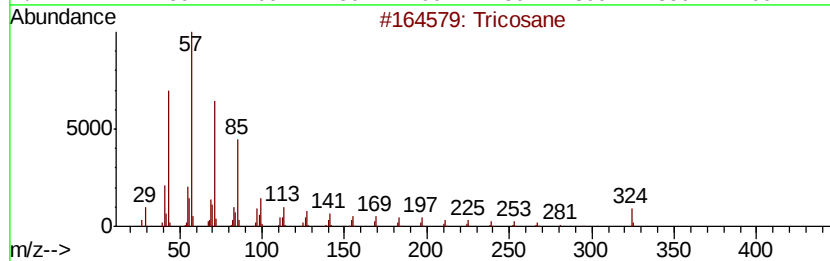
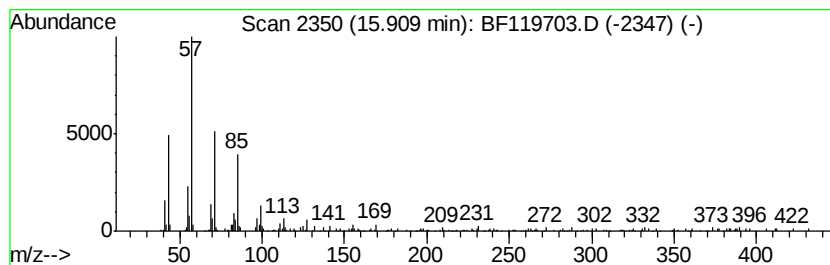
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tricosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	3.83 ng	208727	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	74
2		Octadecane	254	C18H38	000593-45-3	72
3		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	72
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
5		Eicosane	282	C20H42	000112-95-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040120\
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 Operator : MA/SJ
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 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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...	5.03	3.3	ng	261335	1	6.82	1572840	20.0
n-Hexadecanoic acid	11.87	5.1	ng	452127	4	11.34	1762800	20.0
9,12-Octadecadien...	12.42	4.7	ng	416118	4	11.34	1762800	20.0
16-Octadecenoic a...	12.45	6.6	ng	579123	4	11.34	1762800	20.0
Octadecanoic acid	12.66	2.6	ng	225559	4	11.34	1762800	20.0
Tricosyl heptaflu...	13.83	8.5	ng	531773	5	13.98	1251930	20.0
Heptadecane, 2,6,...	14.46	6.4	ng	402244	5	13.98	1251930	20.0
Octadecane	14.76	3.8	ng	207854	6	15.44	1091100	20.0
Nonadecane	15.10	4.7	ng	258285	6	15.44	1091100	20.0
Tricosane	15.91	3.8	ng	208727	6	15.44	1091100	20.0