

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012920\
 Data File : BF118727.D
 Acq On : 29 Jan 2020 13:19
 Operator : CG/JU
 Sample : L1289-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-26-27-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	2.157	5	12	19	rVB	747490	1015196	13.51%	1.553%
2	5.475	570	576	579	rBV	5039124	5099816	67.85%	7.802%
3	6.475	741	746	751	rVB	4843587	5342734	71.08%	8.174%
4	6.845	805	809	813	rBV	1962834	1709476	22.74%	2.615%
5	7.004	832	836	840	rVB	4998919	4518328	60.11%	6.912%
6	7.404	899	904	907	rBV	3481772	3506932	46.66%	5.365%
7	8.128	1023	1027	1031	rBV	2483133	2237927	29.77%	3.424%
8	9.204	1205	1210	1213	rBV	7153747	6850599	91.14%	10.481%
9	9.322	1222	1230	1235	rBV2	156748	167350	2.23%	0.256%
10	9.592	1272	1276	1280	rBV	1251321	1050437	13.98%	1.607%
11	9.880	1320	1325	1329	rBV	2859969	2788490	37.10%	4.266%
12	10.669	1453	1459	1462	rBV	5184652	4846990	64.48%	7.415%
13	11.369	1573	1578	1581	rBV	3068392	2966553	39.47%	4.538%
14	11.592	1612	1616	1626	rBV	169687	186684	2.48%	0.286%
15	11.892	1658	1667	1671	rBV	646762	663951	8.83%	1.016%
16	12.145	1707	1710	1713	rBV	106404	91911	1.22%	0.141%
17	12.598	1776	1787	1794	rBV6	47472	147072	1.96%	0.225%
18	12.674	1797	1800	1812	rVV2	240725	298181	3.97%	0.456%
19	12.951	1842	1847	1850	rBV	7771717	7516533	100.00%	11.499%
20	13.527	1938	1945	1952	rVB2	92046	111735	1.49%	0.171%
21	13.845	1995	1999	2010	rBV	842944	1014189	13.49%	1.552%
22	14.004	2018	2026	2029	rBV	2650858	2665375	35.46%	4.078%
23	14.168	2050	2054	2058	rVB	390622	368494	4.90%	0.564%
24	14.474	2101	2106	2119	rVB	840106	881939	11.73%	1.349%
25	14.780	2150	2158	2163	rBV	1150779	1402095	18.65%	2.145%
26	14.857	2167	2171	2177	rVB	466706	452984	6.03%	0.693%
27	15.115	2209	2215	2225	rVB	1230173	1567739	20.86%	2.398%
28	15.462	2268	2274	2276	rBV	1650953	2144336	28.53%	3.281%
29	15.492	2276	2279	2288	rVB	1071482	1483430	19.74%	2.269%
30	15.927	2346	2353	2359	rBV	727046	1177937	15.67%	1.802%
31	15.974	2359	2361	2371	rVB2	56616	110704	1.47%	0.169%
32	16.433	2432	2439	2450	rVB	337400	637213	8.48%	0.975%
33	17.021	2533	2539	2546	rVB2	85071	178501	2.37%	0.273%
34	17.109	2548	2554	2562	rBV2	30531	76284	1.01%	0.117%

LSC Area Percent Report

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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-26-27-COMP

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

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

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

35	17.509	2617	2622	2632	rvB2	33750	86492	1.15%	0.132%
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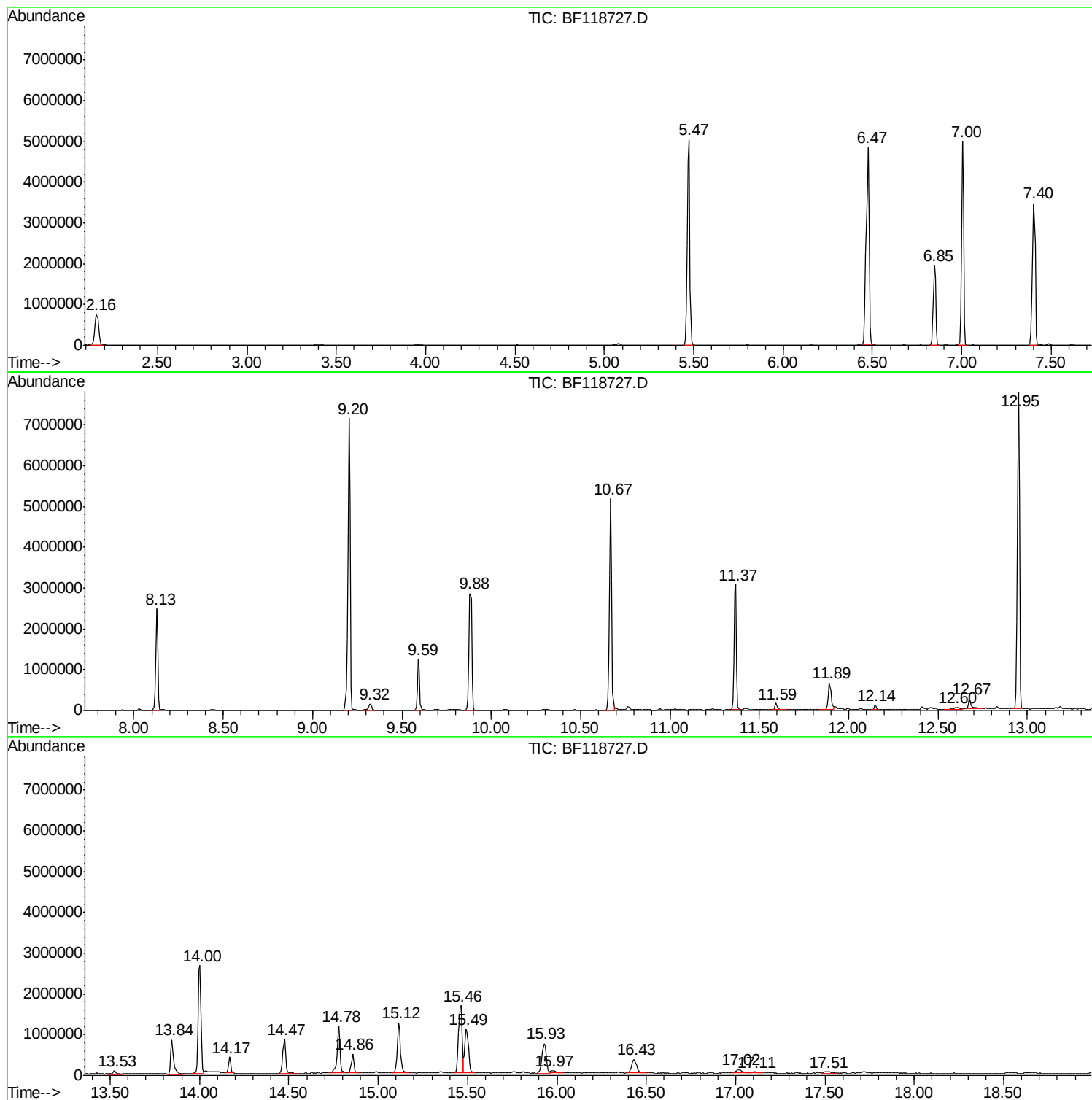
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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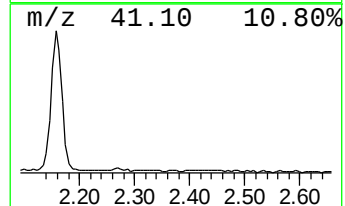
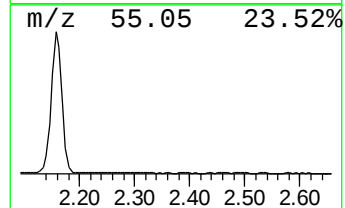
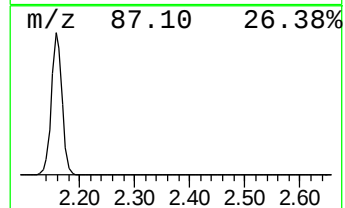
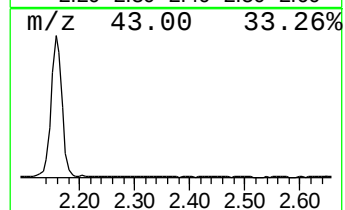
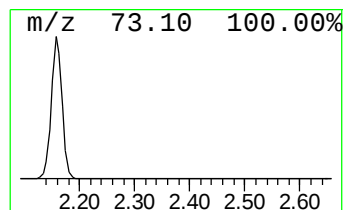
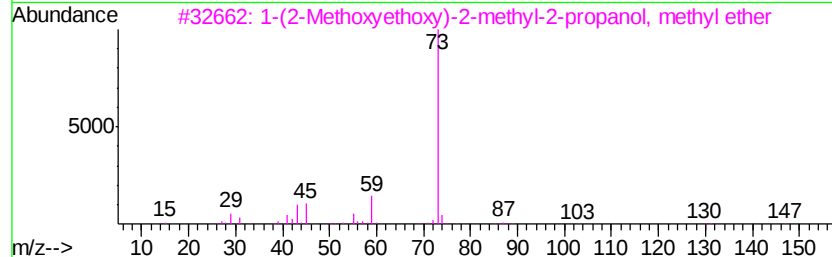
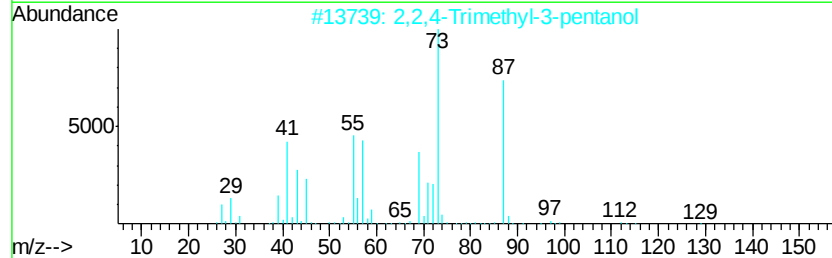
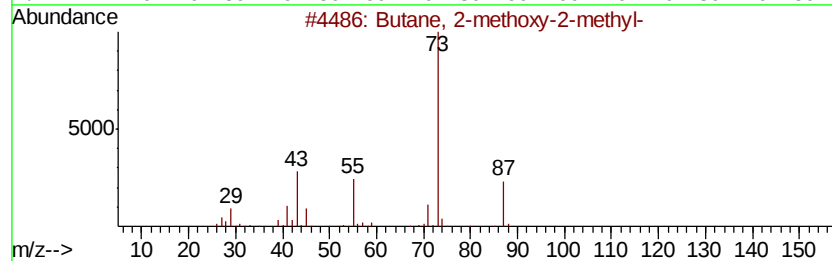
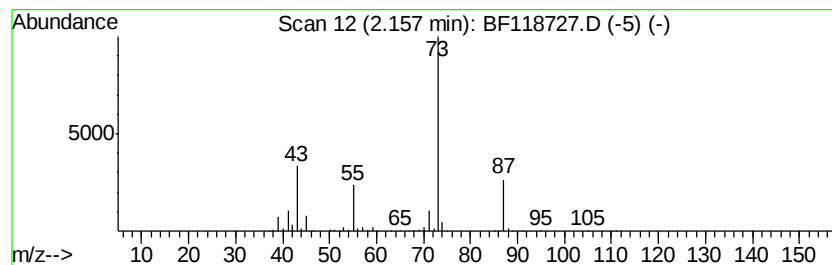
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	11.88 ng	1015200	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3			1-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	23
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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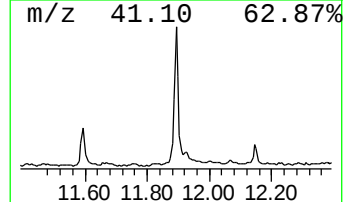
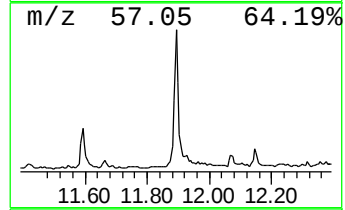
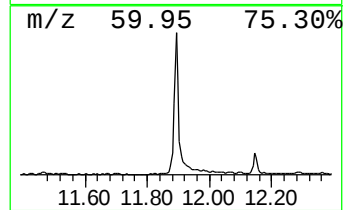
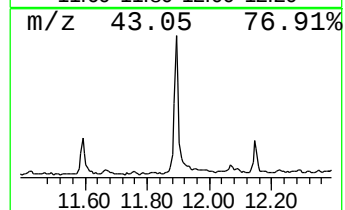
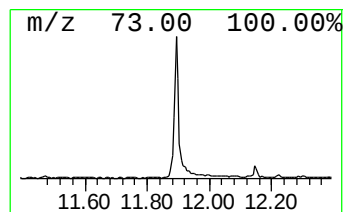
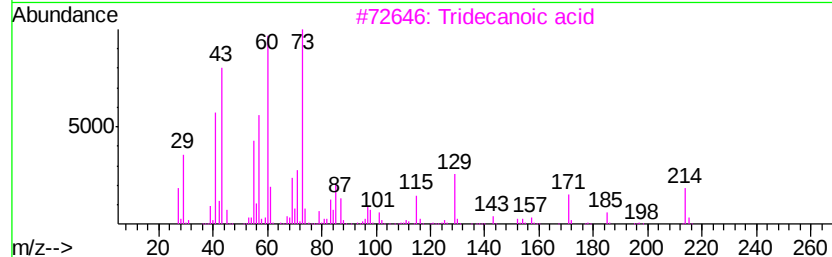
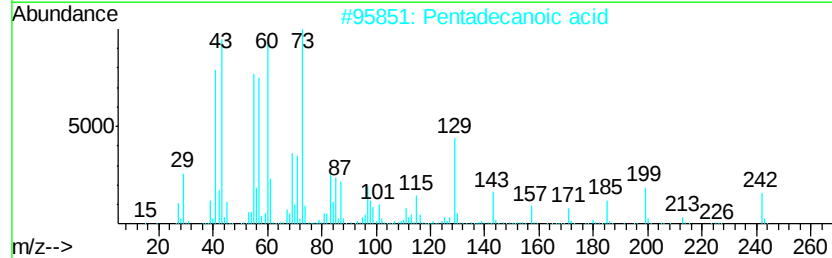
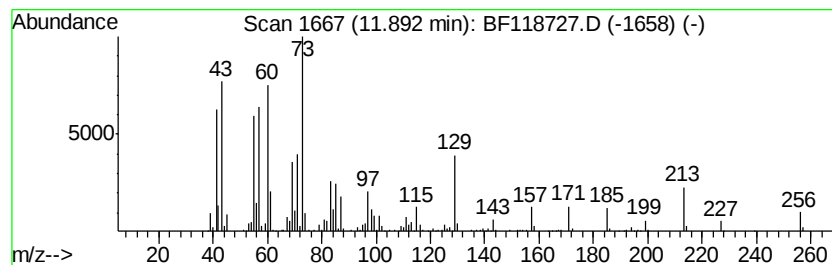
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.89	4.48 ng	663951	Phenanthrene-d10		11.37	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tridecanoic acid	214	C13H26O2	000638-53-9	74
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	70



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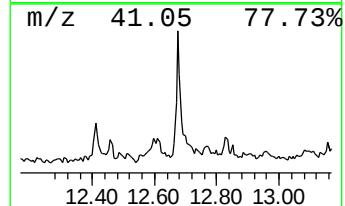
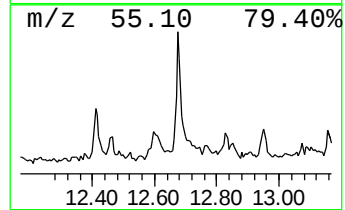
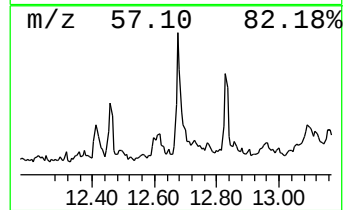
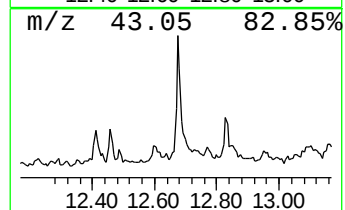
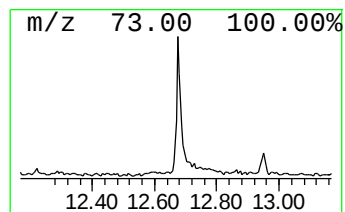
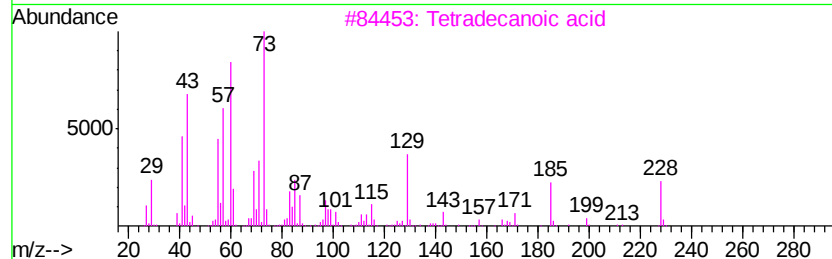
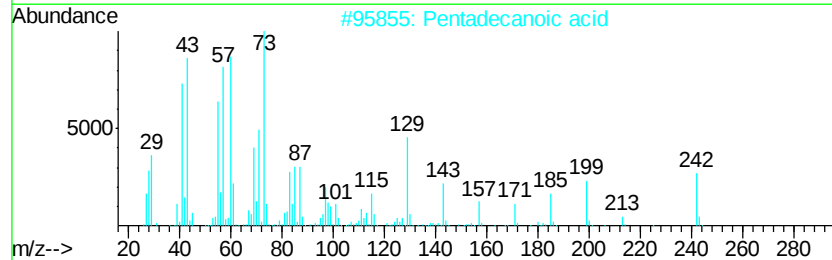
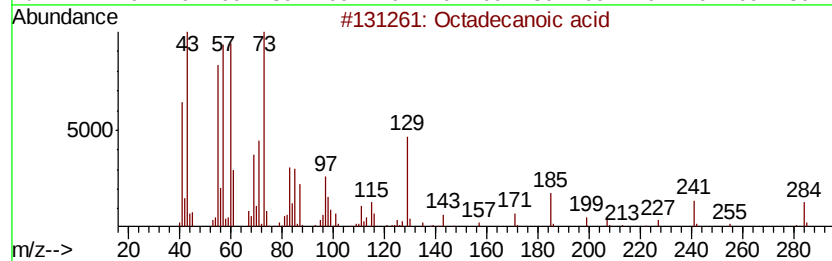
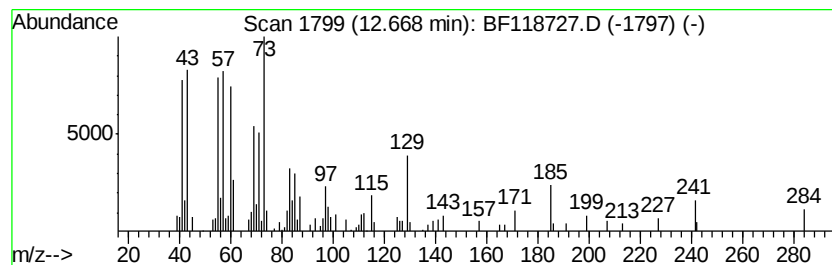
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	2.01 ng	298181	Phenanthrene-d10	11.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4	n-Decanoic acid	172	C10H20O2	000334-48-5	53
5	Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	42



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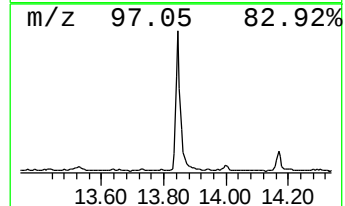
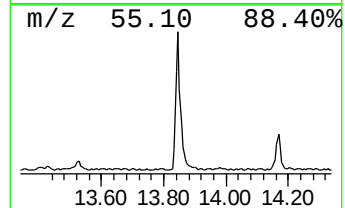
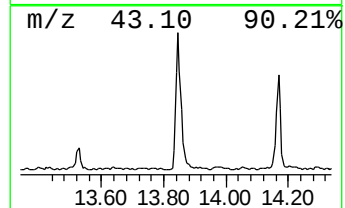
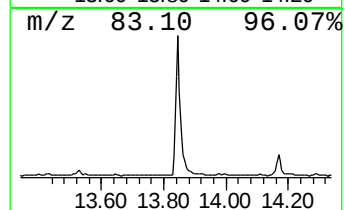
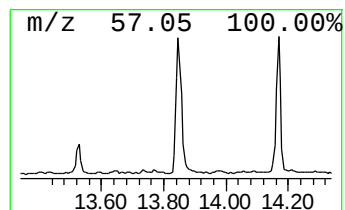
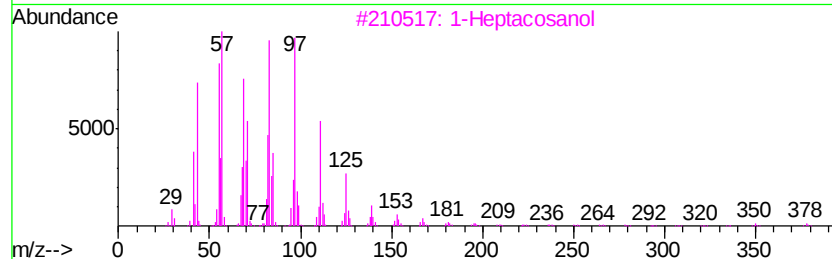
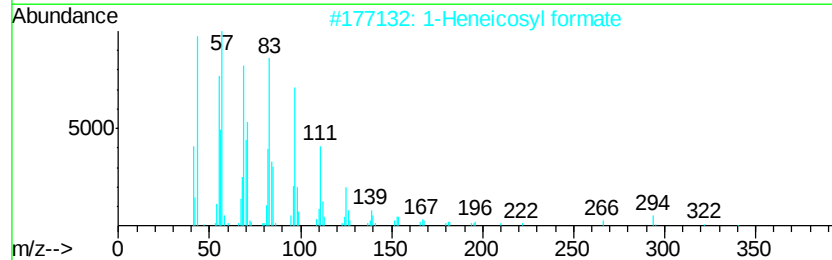
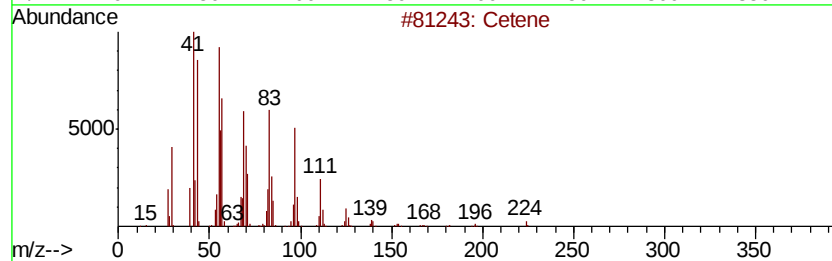
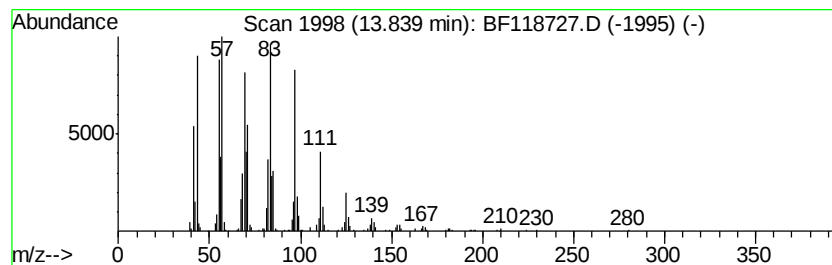
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Cetene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	7.61 ng	1014190	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cetene	224	C16H32	000629-73-2	95
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3		1-Heptacosanol	396	C27H56O	002004-39-9	94
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



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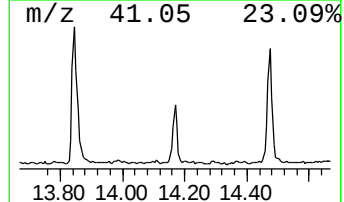
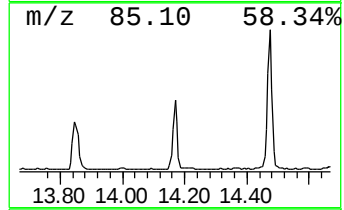
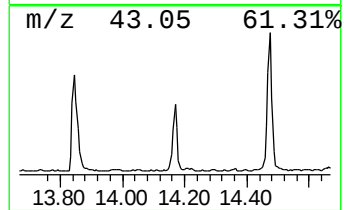
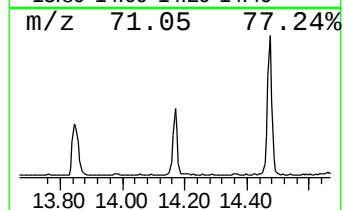
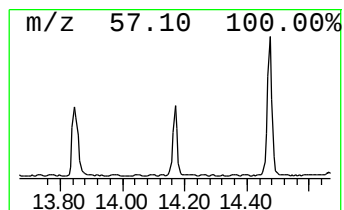
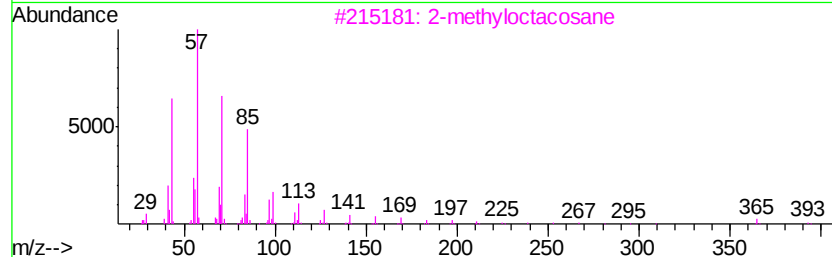
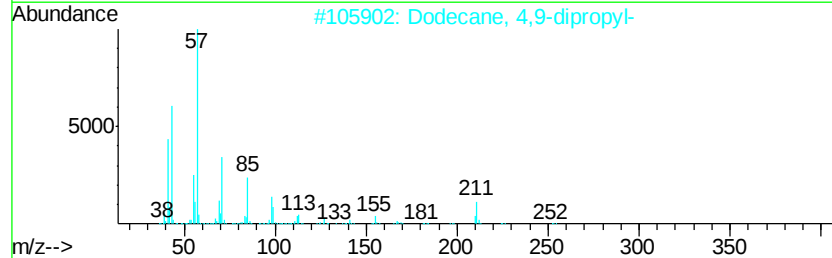
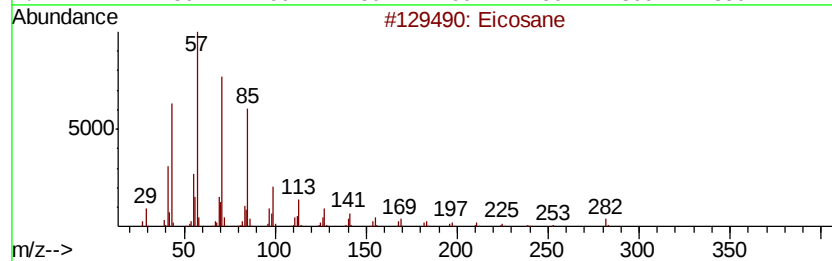
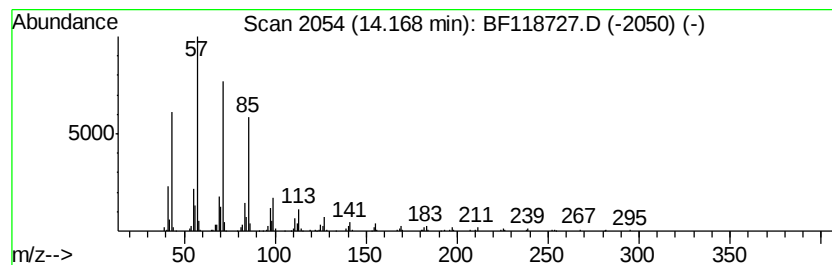
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	2.77 ng	368494	Chrysene-d12	14.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	92
2	Dodecane, 4,9-dipropyl-		254	C18H38	003054-63-5	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	Heneicosane		296	C21H44	000629-94-7	91



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ClientSampleId :
TP-26-27-COMP

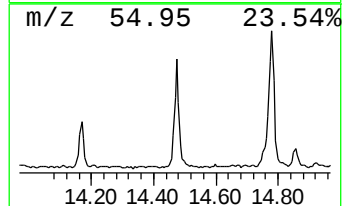
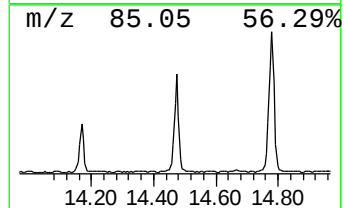
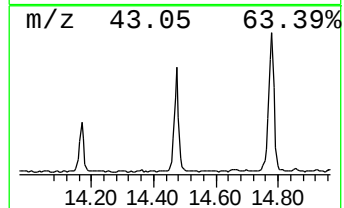
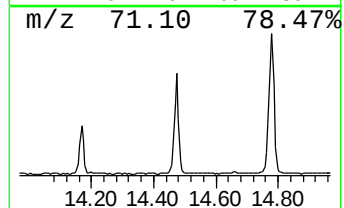
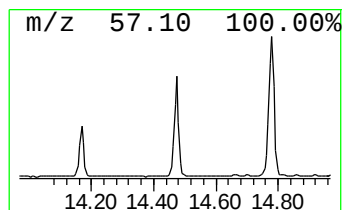
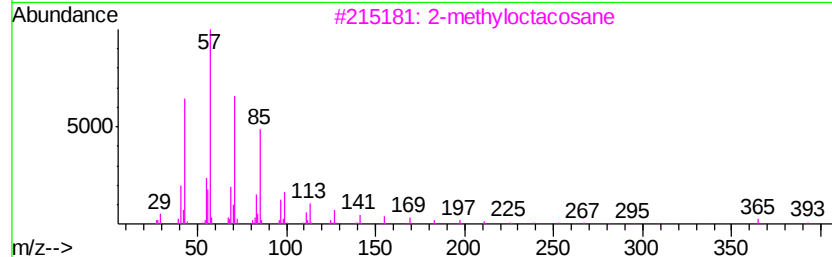
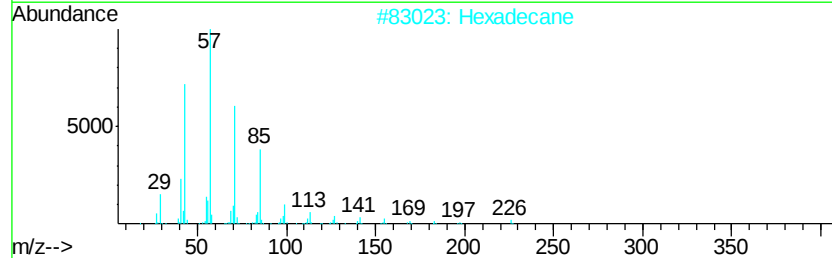
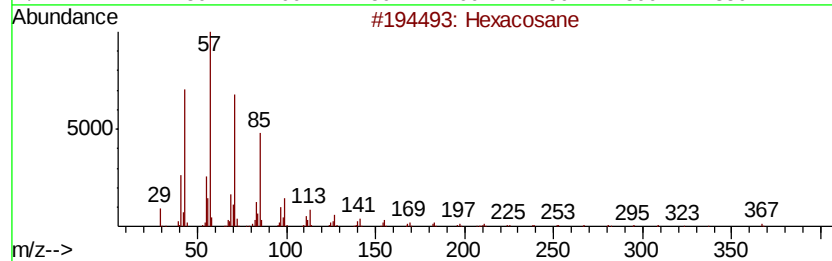
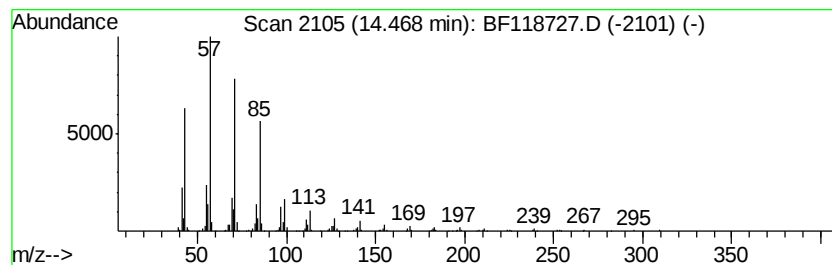
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	6.62 ng	881939	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	94
2		Hexadecane	226	C16H34	000544-76-3	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012920\
 Data File : BF118727.D
 Acq On : 29 Jan 2020 13:19
 Operator : CG/JU
 Sample : L1289-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 TP-26-27-COMP

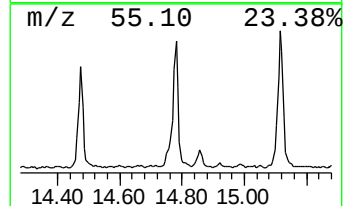
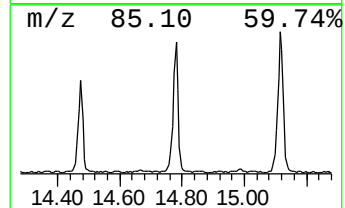
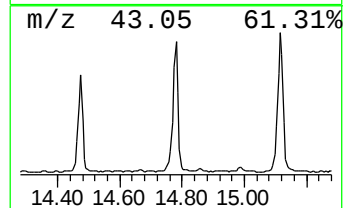
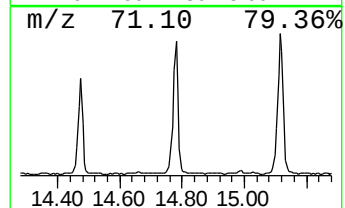
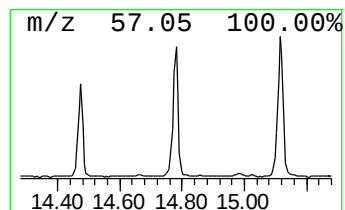
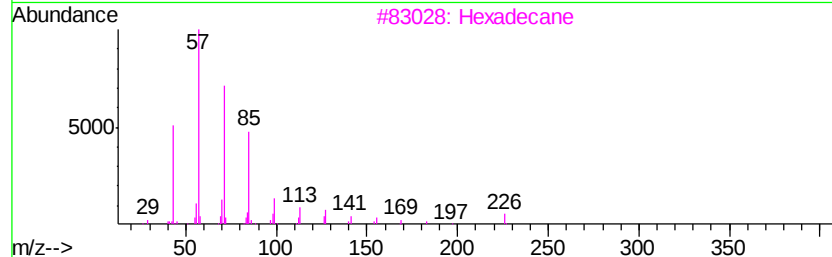
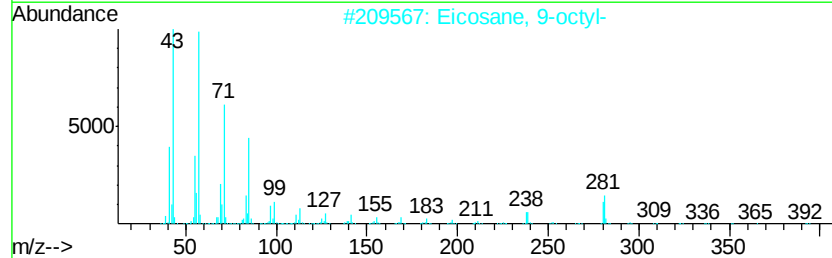
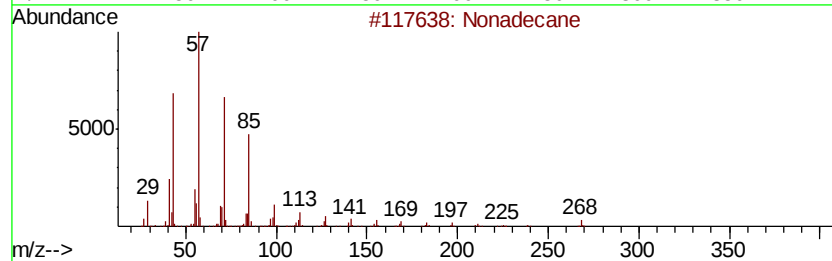
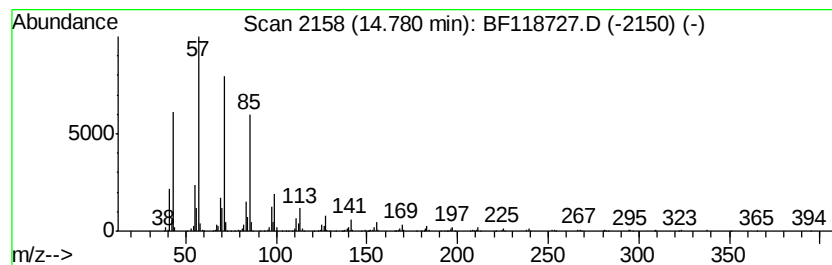
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Nonadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	13.08 ng	1402100	Perylene-d12	15.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	94
2	Eicosane, 9-octyl-		394	C28H58	013475-77-9	93
3	Hexadecane		226	C16H34	000544-76-3	93
4	Hexadecane, 5-butyl-		282	C20H42	006912-07-8	93
5	Heptacosane		380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012920\
 Data File : BF118727.D
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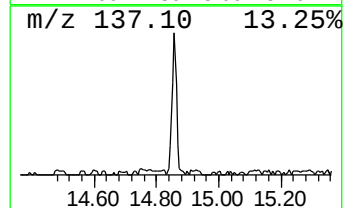
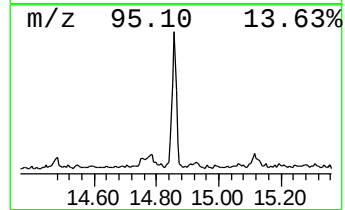
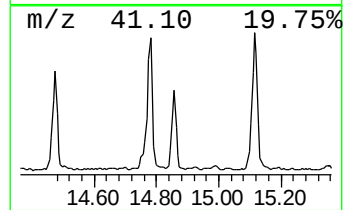
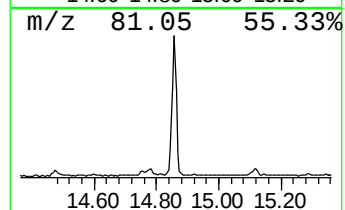
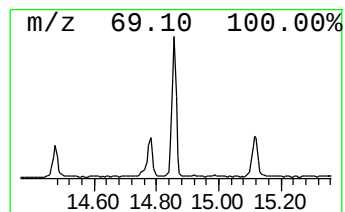
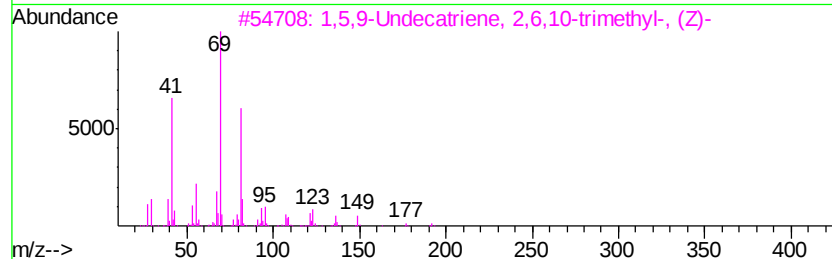
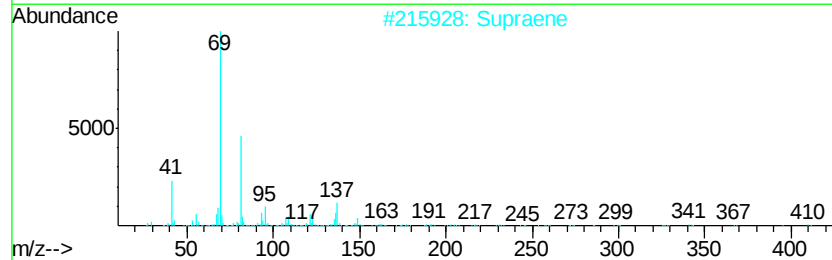
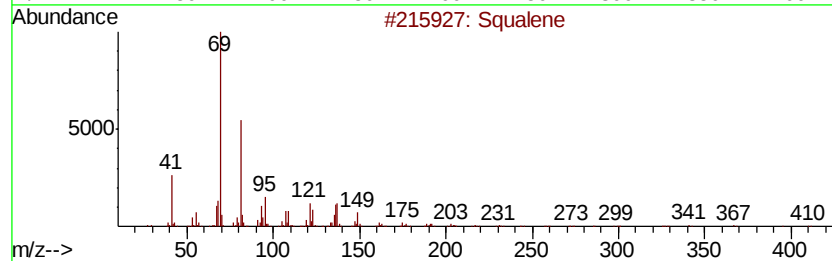
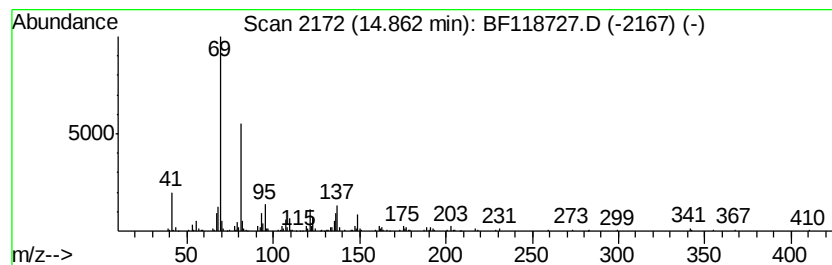
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Squalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.86	4.22 ng	452984	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	99
2		Supraene	410	C30H50	007683-64-9	98
3		1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	87
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
5		Carbonic acid, methyl ester, [(E...	280	C17H28O3	1000164-38-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012920\
Data File : BF118727.D
Acq On : 29 Jan 2020 13:19
Operator : CG/JU
Sample : L1289-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

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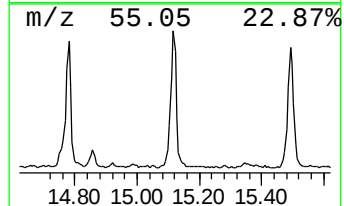
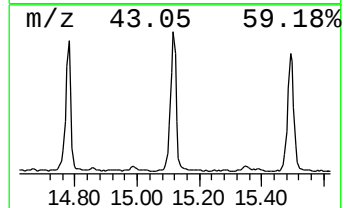
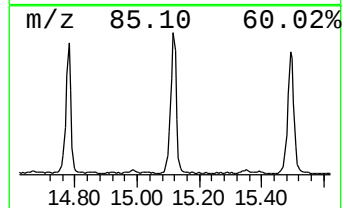
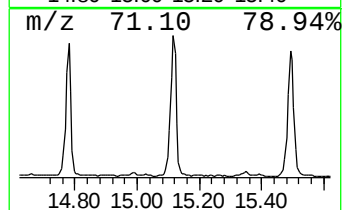
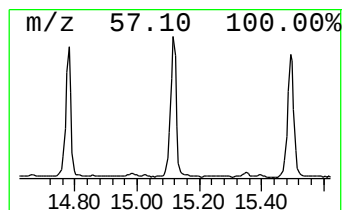
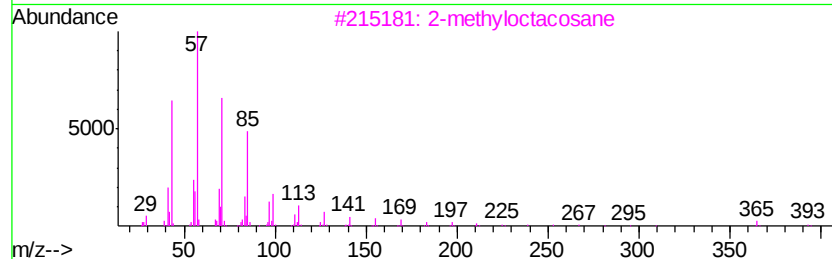
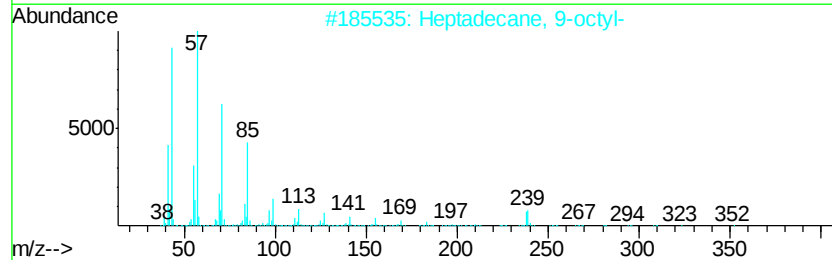
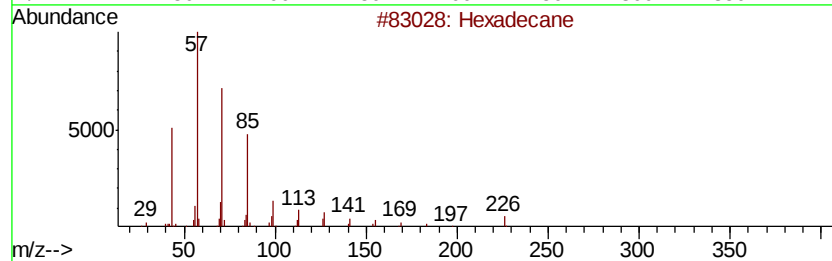
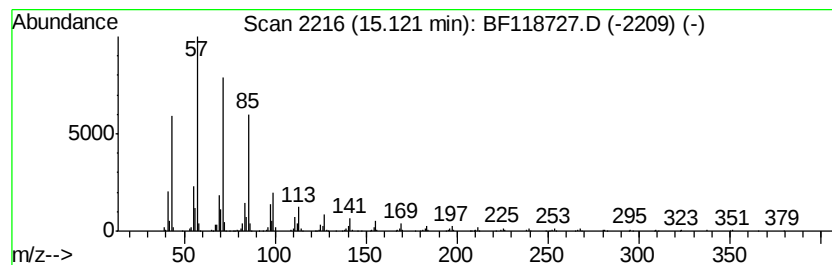
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	14.62 ng	1567740	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hentriacontane	437	C31H64	000630-04-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012920\
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Misc :
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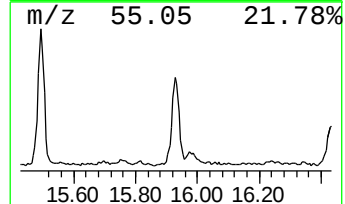
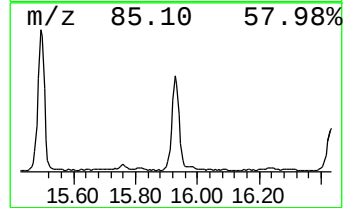
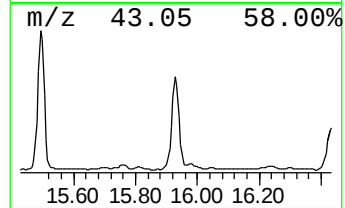
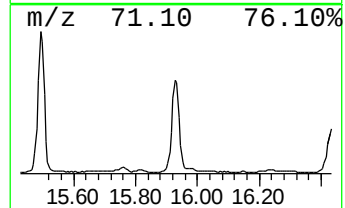
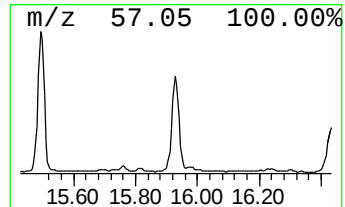
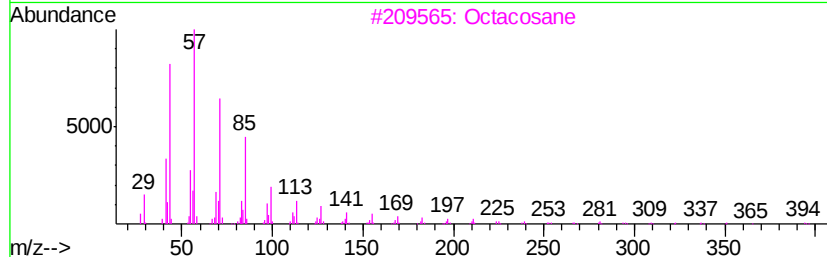
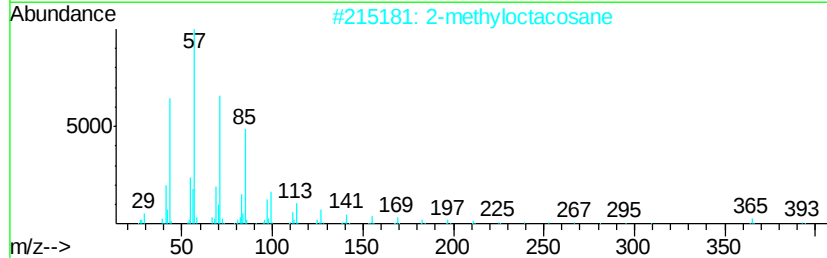
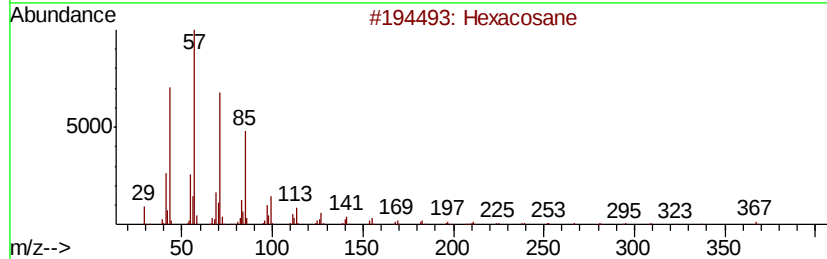
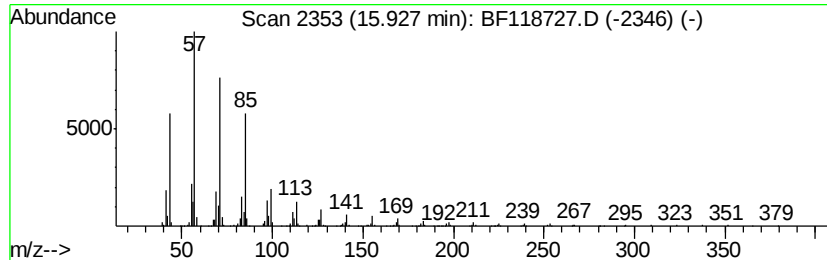
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	10.99 ng	1177940	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012920\
Data File : BF118727.D
Acq On : 29 Jan 2020 13:19
Operator : CG/JU
Sample : L1289-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

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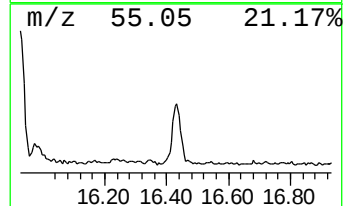
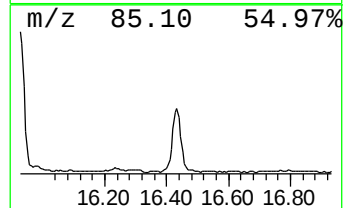
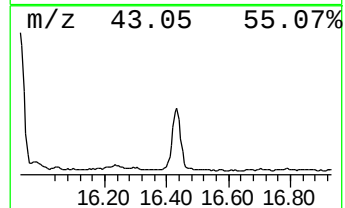
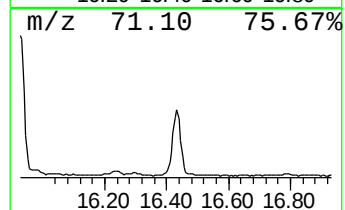
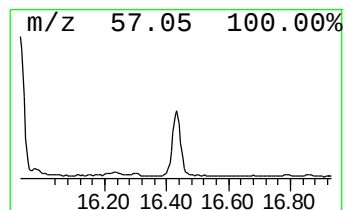
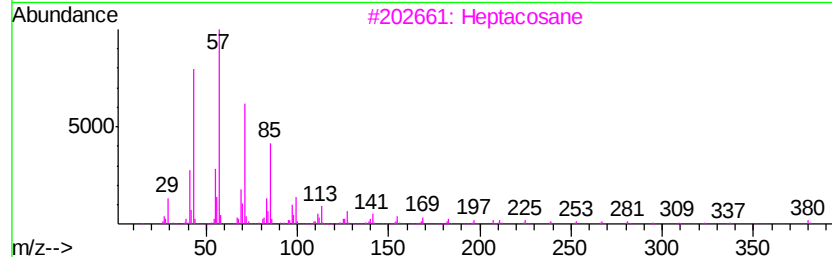
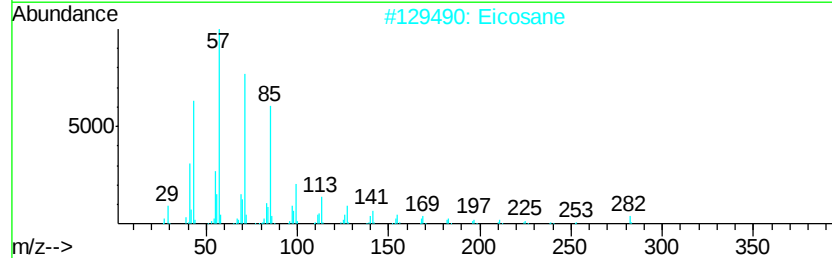
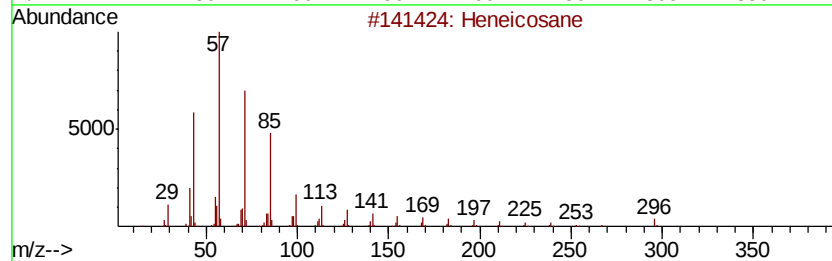
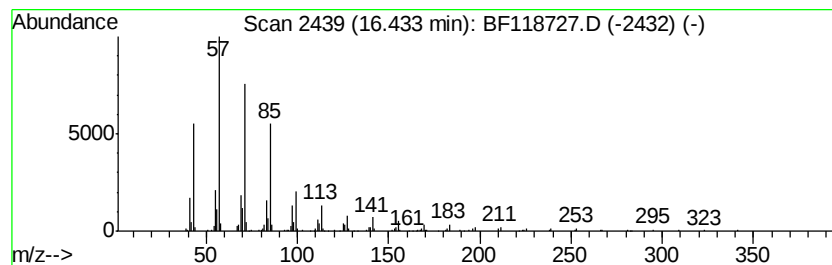
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	5.94 ng	637213	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Triacontane	422	C30H62	000638-68-6	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012920\
 Data File : BF118727.D
 Acq On : 29 Jan 2020 13:19
 Operator : CG/JU
 Sample : L1289-05
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 ALS Vial : 3 Sample Multiplier: 1

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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic acid	11.89	4.5	ng	663951	4	11.37	2966550	20.0
Octadecanoic acid	12.67	2.0	ng	298181	4	11.37	2966550	20.0
Cetene	13.84	7.6	ng	1014190	5	14.00	2665380	20.0
Eicosane	14.17	2.8	ng	368494	5	14.00	2665380	20.0
Hexacosane	14.47	6.6	ng	881939	5	14.00	2665380	20.0
Nonadecane	14.78	13.1	ng	1402100	6	15.46	2144340	20.0
Squalene	14.86	4.2	ng	452984	6	15.46	2144340	20.0
Heptadecane, 9-oc...	15.12	14.6	ng	1567740	6	15.46	2144340	20.0
2-methyloctacosane	15.93	11.0	ng	1177940	6	15.46	2144340	20.0
Heneicosane	16.43	5.9	ng	637213	6	15.46	2144340	20.0