

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
 Data File : BF105081.D
 Acq On : 3 May 2018 19:10
 Operator : JU/SJ
 Sample : J2665-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ES-RT-1871

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:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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	4.998	491	495	510	rBV	50764	92127	3.62%	0.387%
2	5.404	561	564	601	rBV	1542277	2209516	86.77%	9.271%
3	6.434	735	739	756	rBV	1702762	2183583	85.75%	9.163%
4	6.557	756	760	784	rVB	2172313	2369698	93.06%	9.944%
5	6.775	794	797	808	rBV	551118	580597	22.80%	2.436%
6	6.934	820	824	840	rBV	1916698	1813657	71.23%	7.610%
7	7.351	891	895	912	rBV	1341964	1424000	55.92%	5.975%
8	8.063	1013	1016	1034	rBV	795785	829053	32.56%	3.479%
9	9.139	1195	1199	1209	rBV	2757717	2546376	100.00%	10.685%
10	9.539	1264	1267	1274	rVB	227307	229371	9.01%	0.962%
11	9.739	1298	1301	1305	rVB	52187	40491	1.59%	0.170%
12	9.822	1311	1315	1327	rVB	923326	939122	36.88%	3.941%
13	10.239	1383	1386	1389	rVB	81570	62734	2.46%	0.263%
14	10.457	1419	1423	1426	rBV	20679	27694	1.09%	0.116%
15	10.616	1446	1450	1460	rBV	1564455	1556033	61.11%	6.529%
16	10.710	1464	1466	1470	rBV	66313	72562	2.85%	0.304%
17	10.822	1482	1485	1489	rVB3	23250	26760	1.05%	0.112%
18	11.163	1540	1543	1545	rBV	35055	29665	1.16%	0.124%
19	11.310	1565	1568	1574	rBV	862610	875994	34.40%	3.676%
20	11.839	1653	1658	1671	rBV	392990	456458	17.93%	1.915%
21	12.533	1772	1776	1786	rBV	63075	113847	4.47%	0.478%
22	12.621	1787	1791	1800	rVV	123580	135808	5.33%	0.570%
23	12.745	1809	1812	1813	rBV2	32688	28206	1.11%	0.118%
24	12.763	1813	1815	1822	rVB	39374	50165	1.97%	0.210%
25	12.904	1835	1839	1842	rBV	2042727	1874430	73.61%	7.865%
26	13.468	1932	1935	1944	rBV4	26729	42032	1.65%	0.176%
27	13.833	1994	1997	2008	rVB	135256	195764	7.69%	0.821%
28	14.021	2025	2029	2038	rBV	541556	593488	23.31%	2.490%
29	14.115	2042	2045	2050	rVB	88616	81222	3.19%	0.341%
30	14.174	2052	2055	2065	rVB8	31601	66384	2.61%	0.279%
31	14.533	2112	2116	2122	rBV3	74285	121943	4.79%	0.512%
32	14.874	2170	2174	2180	rBV3	70371	112590	4.42%	0.472%
33	15.039	2199	2202	2206	rVB3	36942	38226	1.50%	0.160%
34	15.227	2230	2234	2239	rVB	392823	439613	17.26%	1.845%

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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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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

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35	15.598	2292	2297	2303	rBV	449195	668433	26.25%	2.805%
36	15.815	2330	2334	2338	rBV3	49165	67674	2.66%	0.284%
37	16.045	2368	2373	2381	rBV	285674	432404	16.98%	1.814%
38	16.298	2412	2416	2426	rVB10	85937	177309	6.96%	0.744%
39	16.839	2503	2508	2514	rVB4	62143	114929	4.51%	0.482%
40	17.339	2588	2593	2601	rVB8	47760	111520	4.38%	0.468%

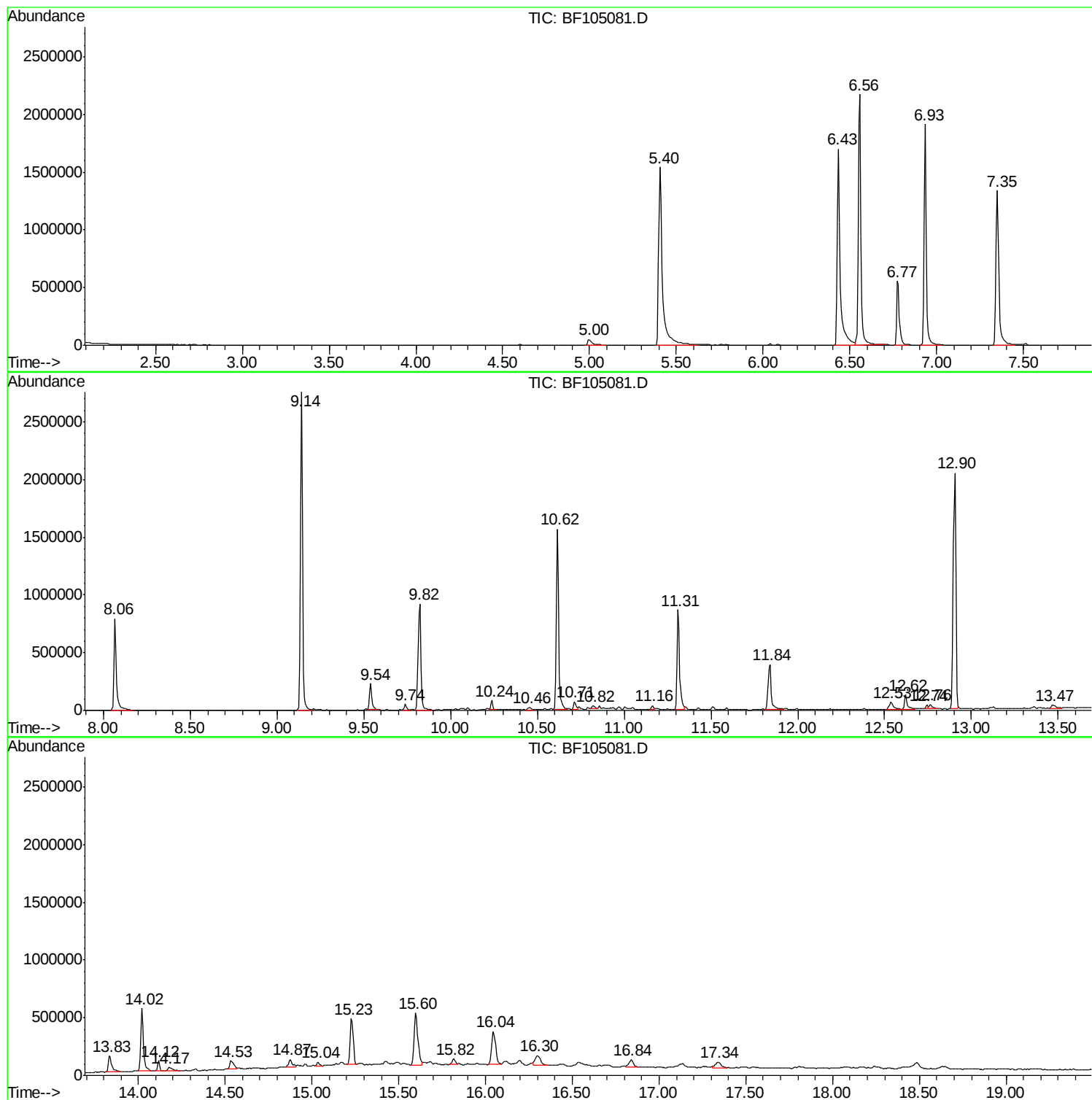
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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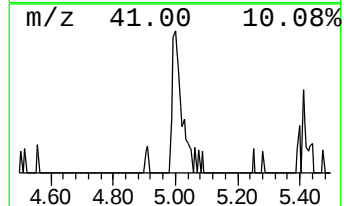
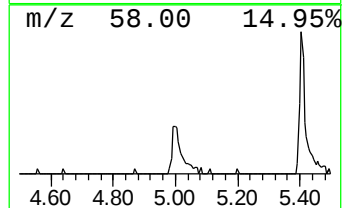
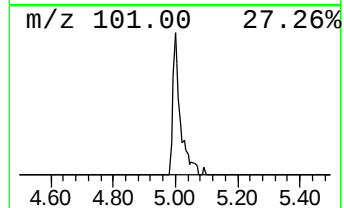
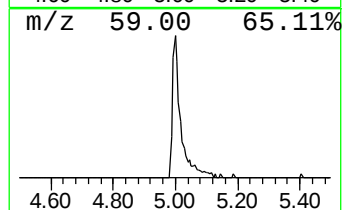
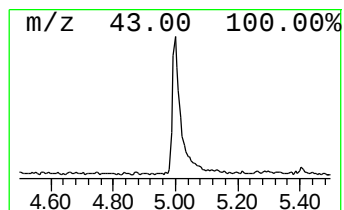
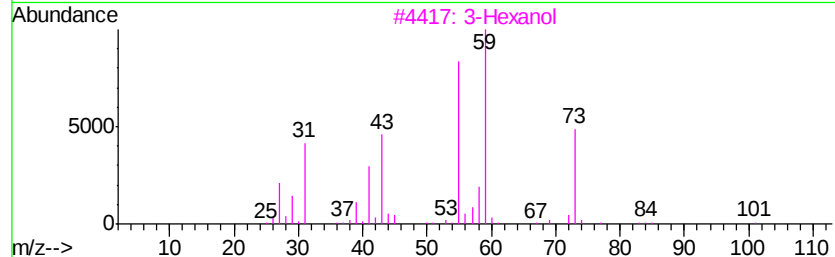
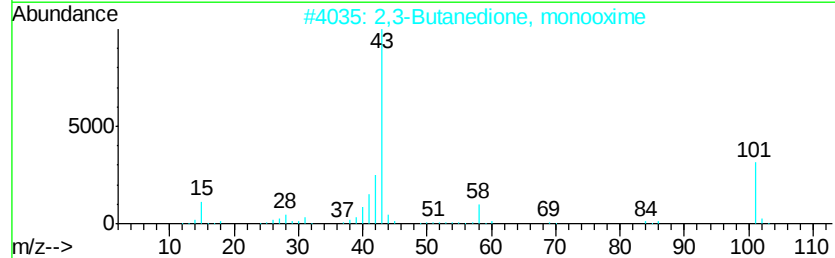
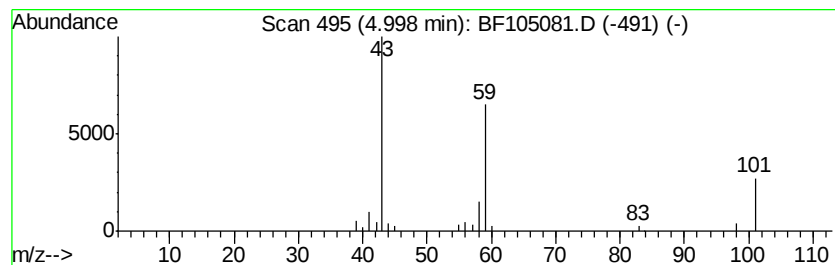
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	3.17 ng	92127	1,4-Dichlorobenzene-d4	6.77

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3		3-Hexanol	102	C6H14O	000623-37-0	28
4		3-Pentanol	88	C5H12O	000584-02-1	25
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25



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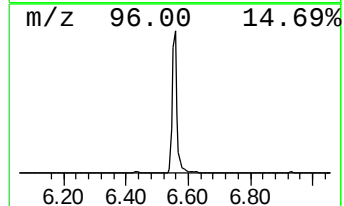
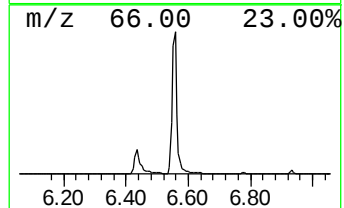
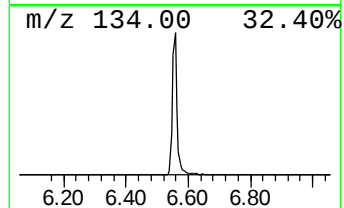
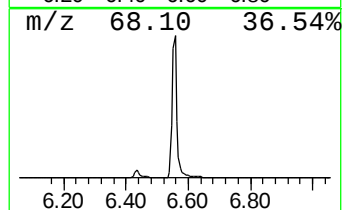
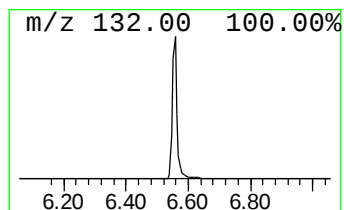
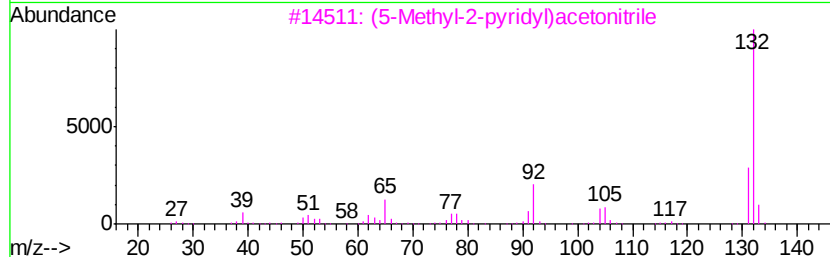
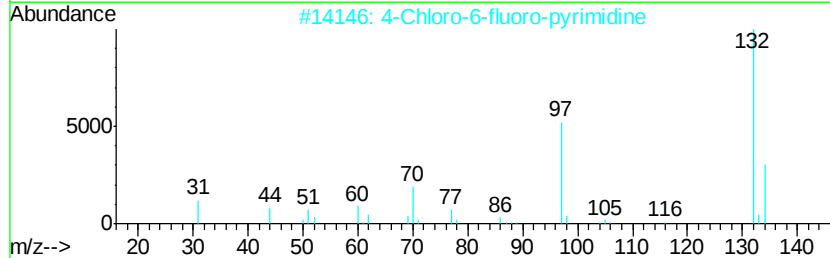
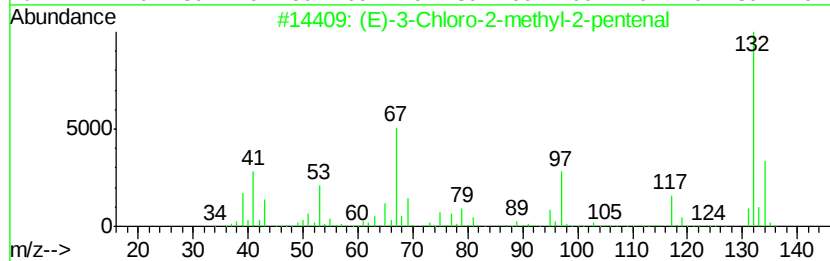
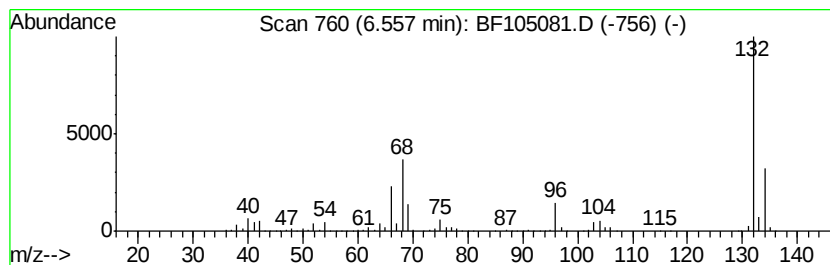
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	81.63 ng	2369700	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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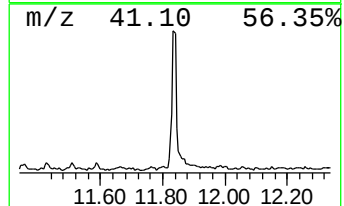
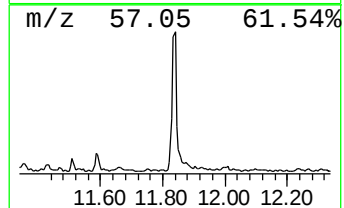
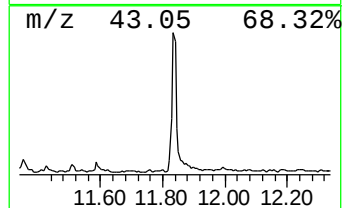
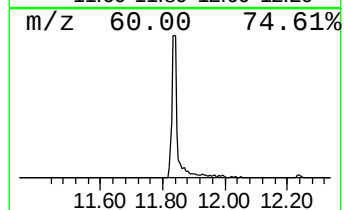
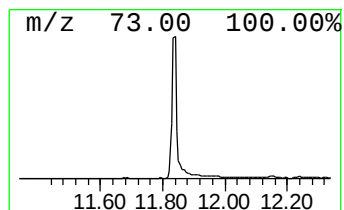
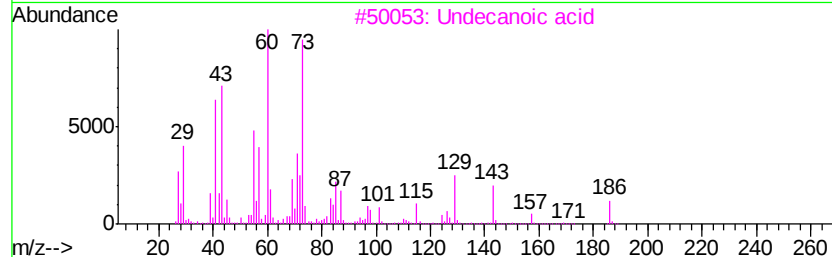
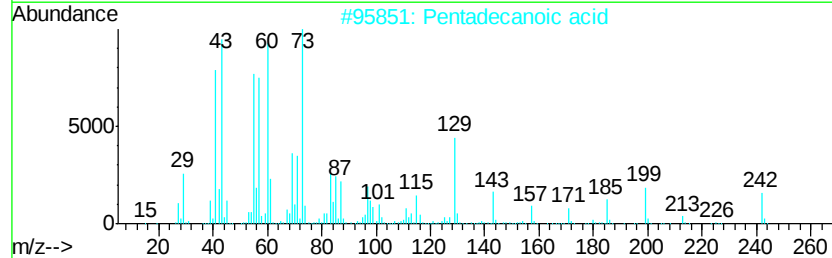
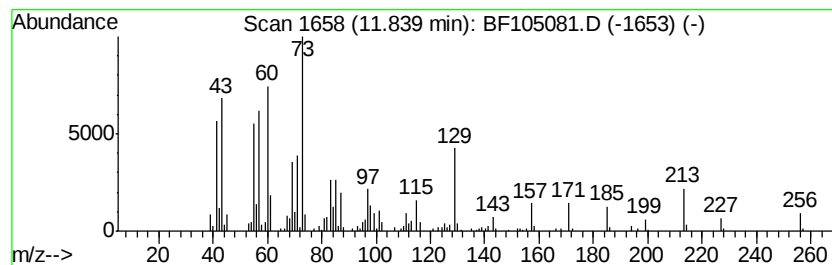
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	10.42 ng	456458	Phenanthrene-d10	11.31

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	87
3		Undecanoic acid	186	C11H22O2	000112-37-8	72
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Tridecanoic acid	214	C13H26O2	000638-53-9	64



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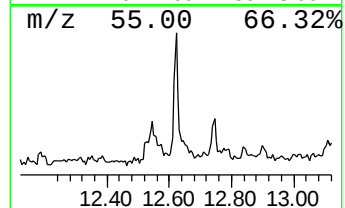
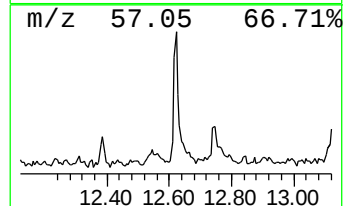
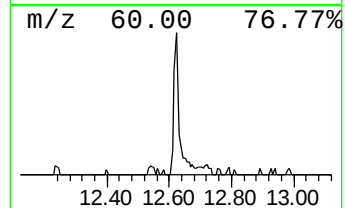
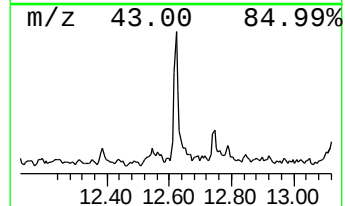
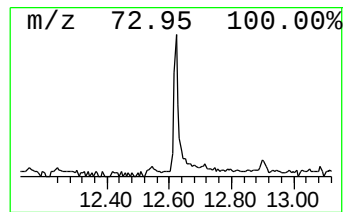
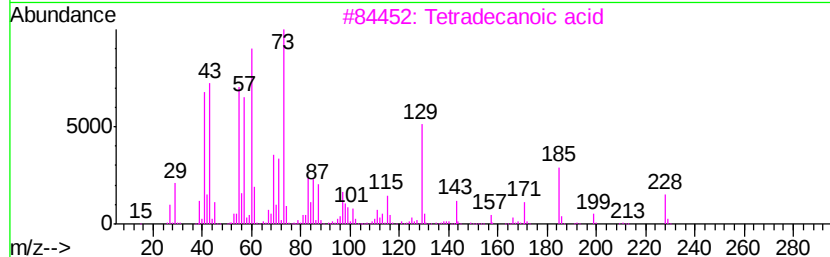
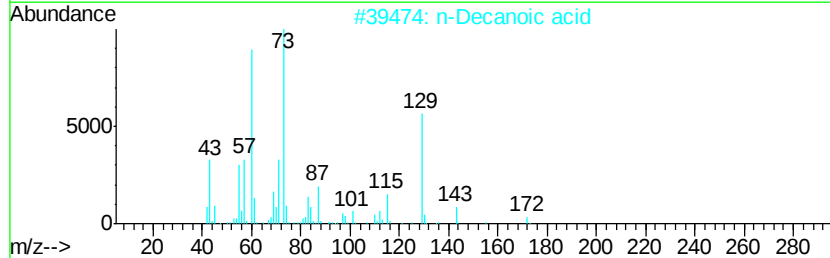
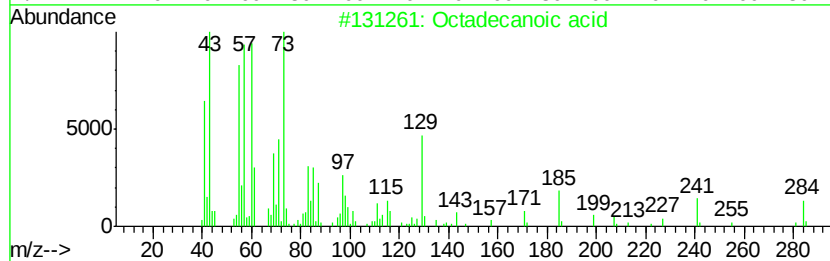
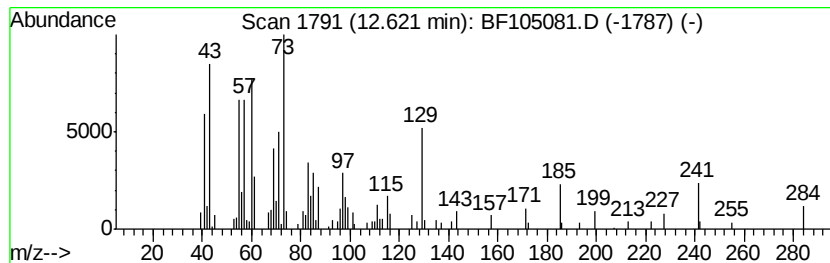
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.62	3.10 ng	135808	Phenanthrene-d10	11.31	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	n-Decanoic acid	172	C10H20O2	000334-48-5	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	41
5	Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	30



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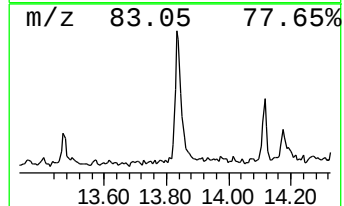
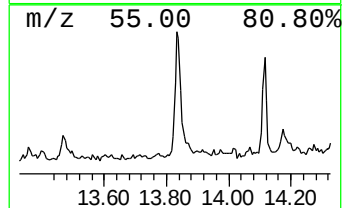
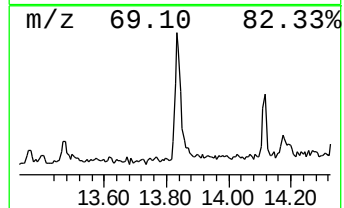
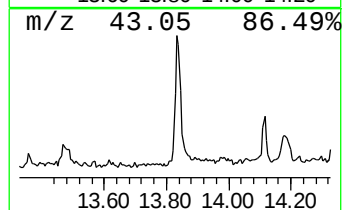
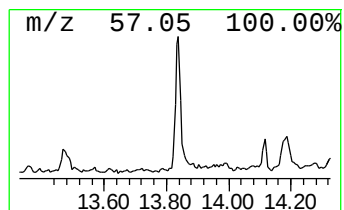
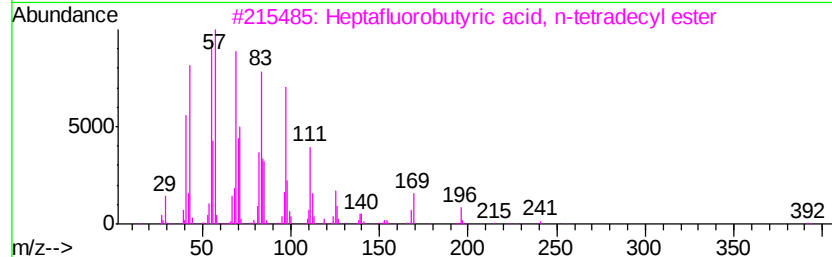
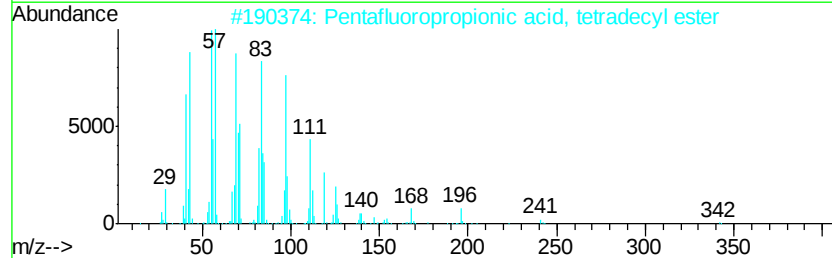
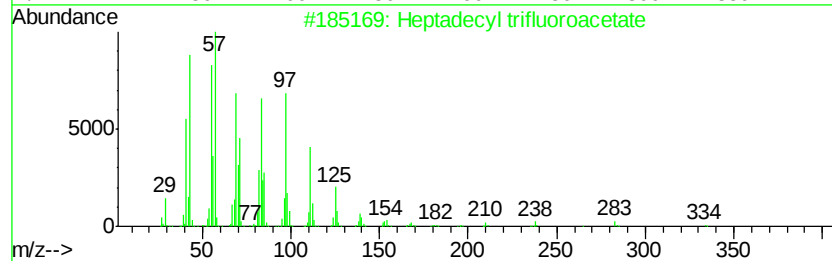
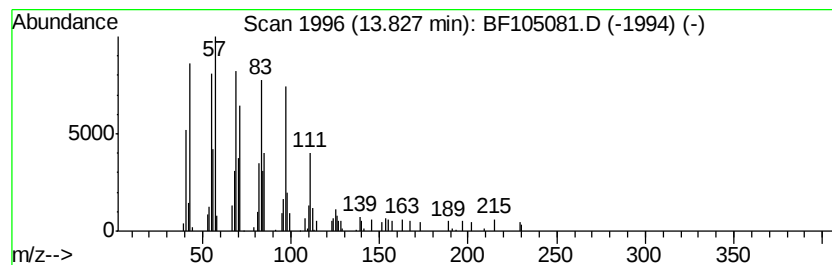
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Heptadecyl trifluoroacetate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	6.60 ng	195764	Chrysene-d12	14.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	95
2			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	94
3			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
4			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94



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Misc :
ALS Vial : 10 Sample Multiplier: 1

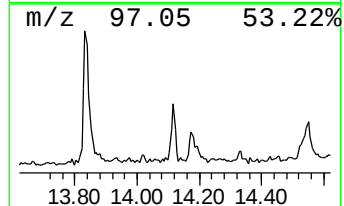
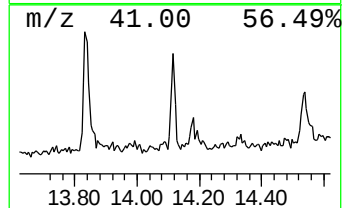
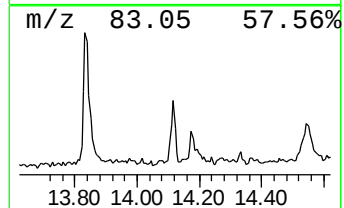
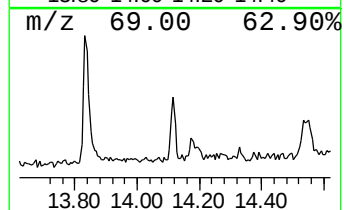
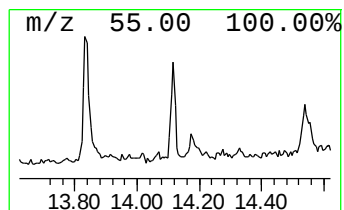
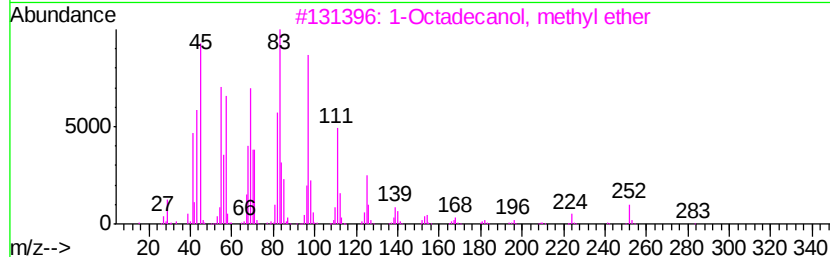
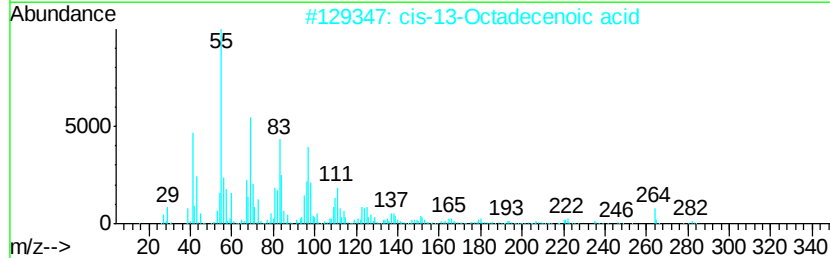
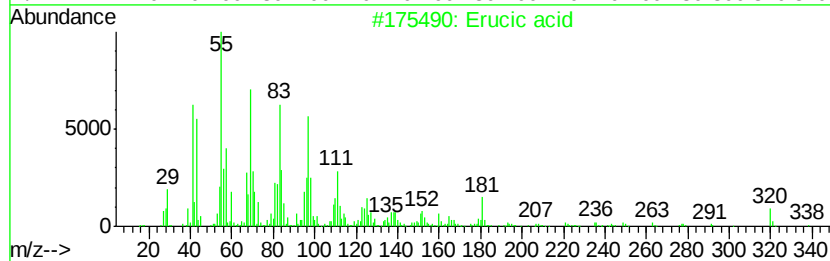
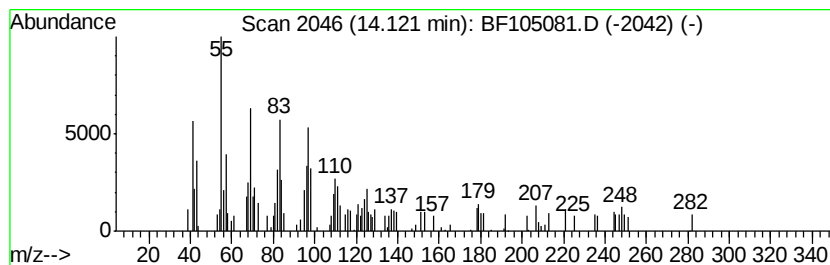
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Erucic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	2.74 ng	81222	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Erucic acid		338 C22H42O2	000112-86-7 87
2	cis-13-Octadecenoic acid		282 C18H34O2	013126-39-1 83
3	1-Octadecanol, methyl ether		284 C19H40O	1000333-92-7 70
4	1-Pentadecanethiol		244 C15H32S	025276-70-4 68
5	1-Tetradecanol		214 C14H30O	000112-72-1 64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
Data File : BF105081.D
Acq On : 3 May 2018 19:10
Operator : JU/SJ
Sample : J2665-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ES-RT-1871

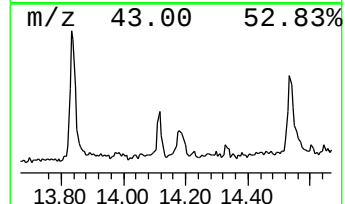
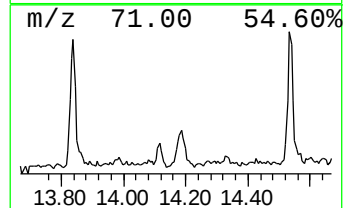
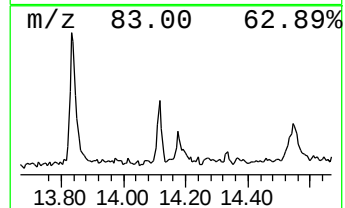
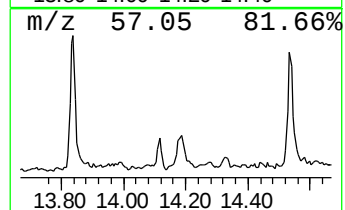
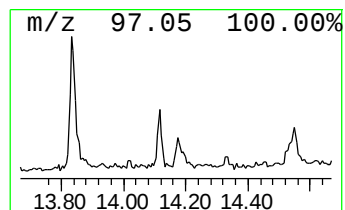
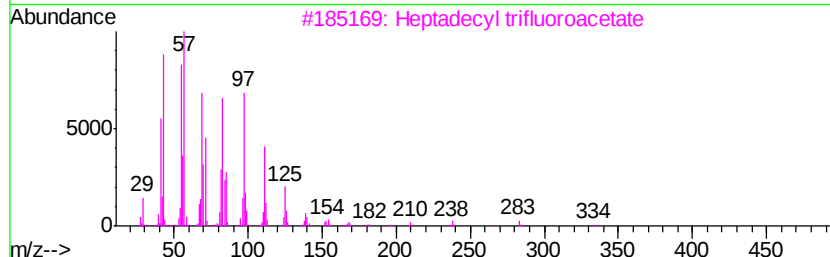
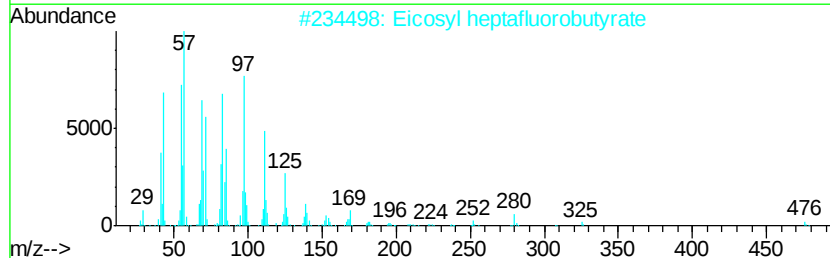
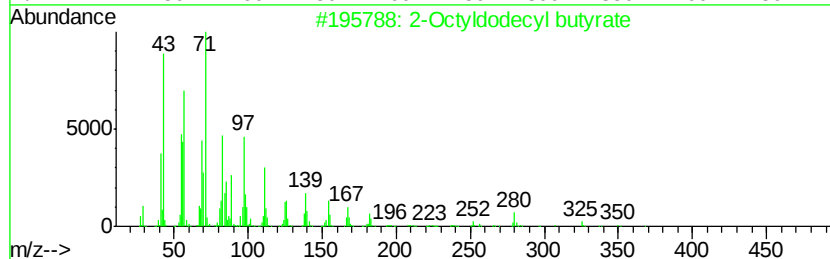
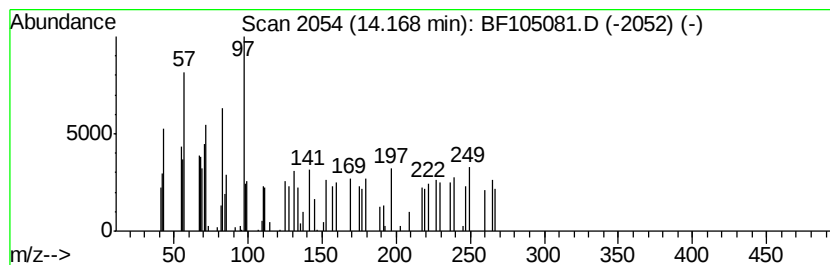
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-Octyldodecyl butyrate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	2.24 ng	66384	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Octyldodecyl butyrate	368	C24H48O2	1000368-55-2	86
2		Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	83
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	83
4		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	83
5		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
Data File : BF105081.D
Acq On : 3 May 2018 19:10
Operator : JU/SJ
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ES-RT-1871

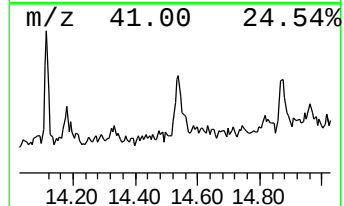
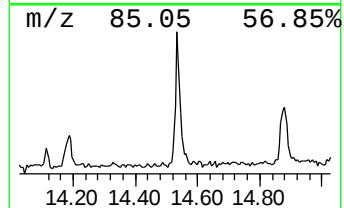
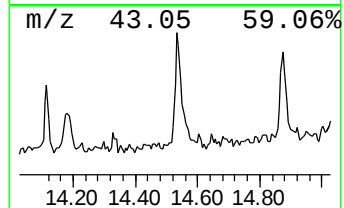
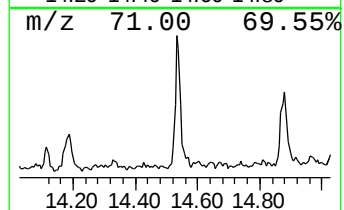
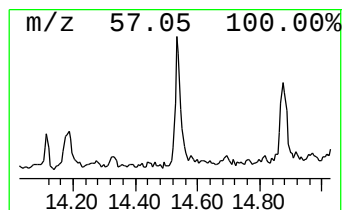
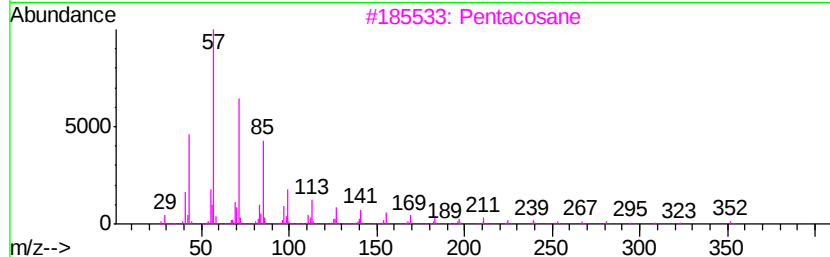
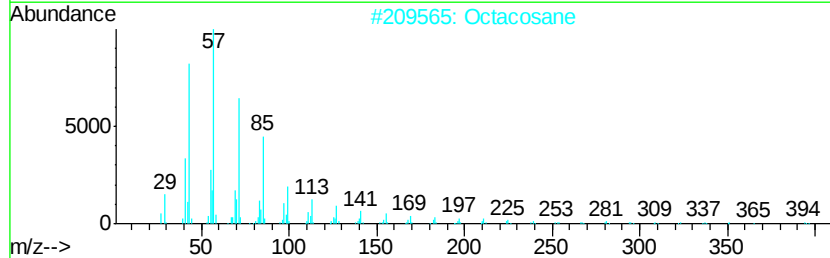
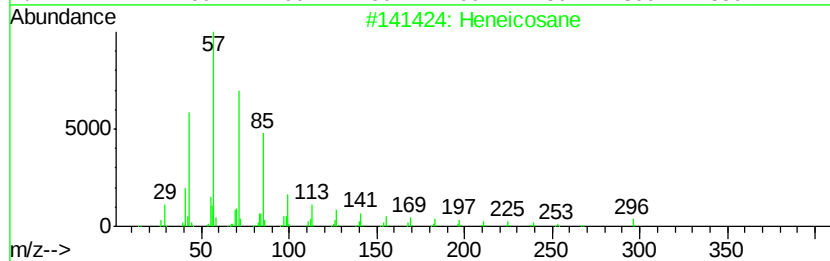
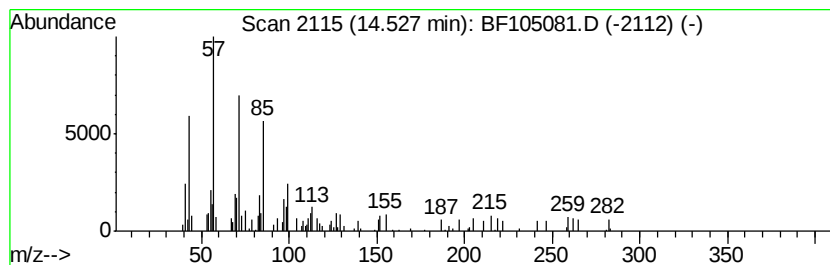
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	4.11 ng	121943	Chrysene-d12	14.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	87
2			Octacosane	394	C28H58	000630-02-4	87
3			Pentacosane	352	C25H52	000629-99-2	87
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
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Acq On : 3 May 2018 19:10
Operator : JU/SJ
Sample : J2665-09
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ALS Vial : 10 Sample Multiplier: 1

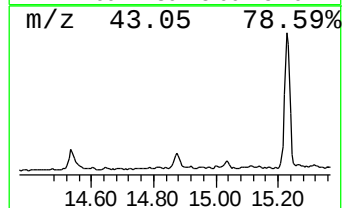
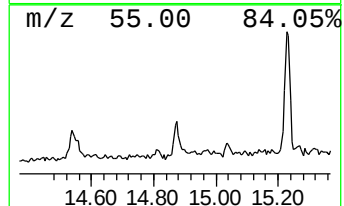
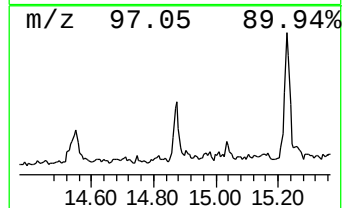
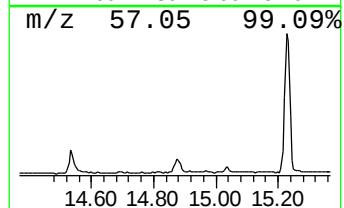
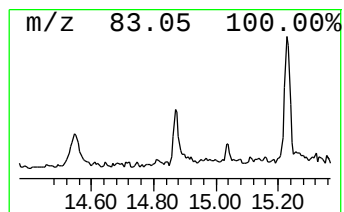
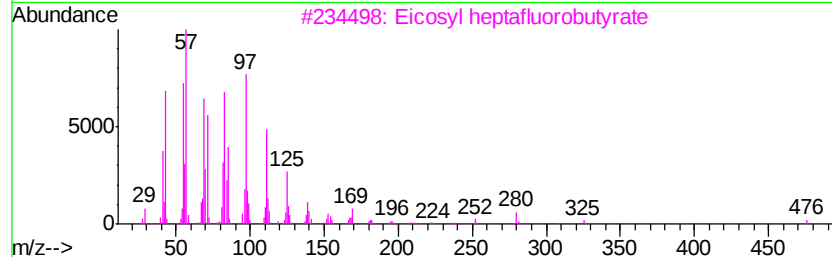
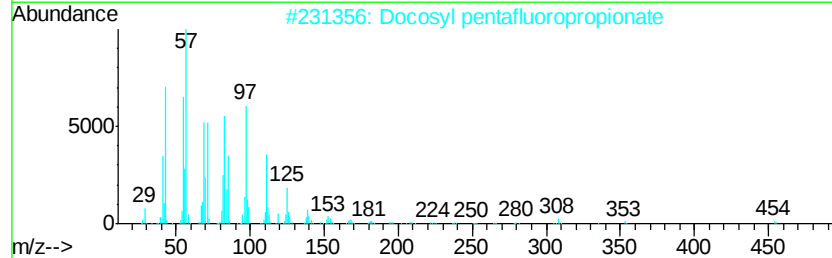
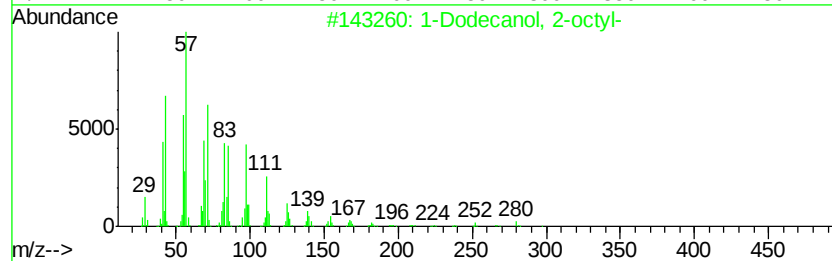
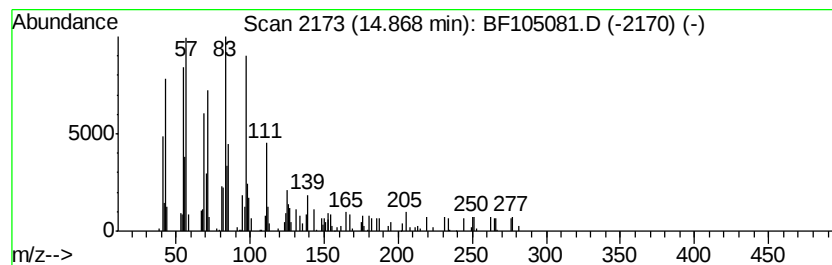
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Dodecanol, 2-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.87	3.37 ng	112590	Perylene-d12	15.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	87
2	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	87
3	Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	83
4	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	83
5	Octatriacontyl trifluoroacetate	647	C40H77F3O2	1000351-88-7	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050318\
Data File : BF105081.D
Acq On : 3 May 2018 19:10
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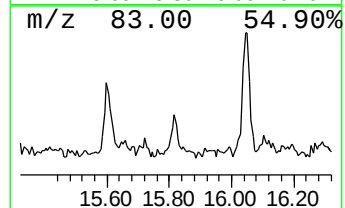
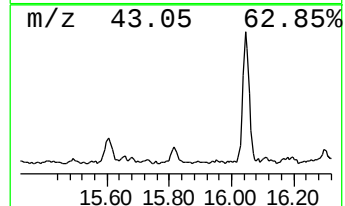
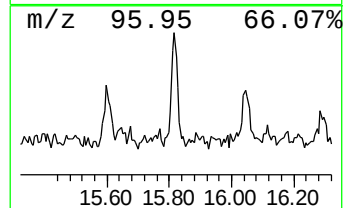
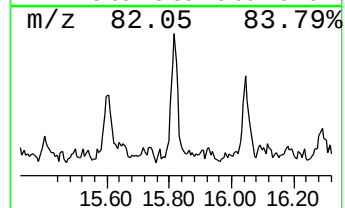
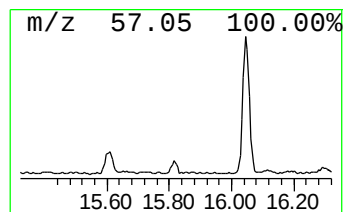
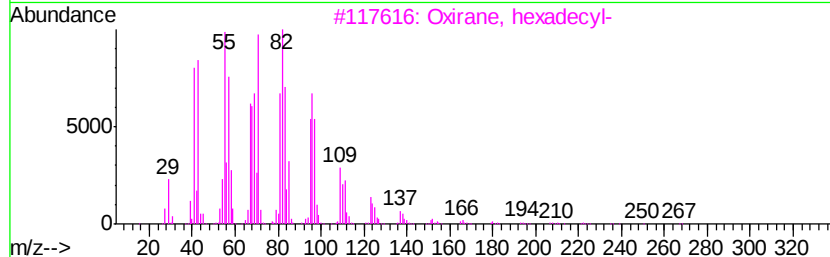
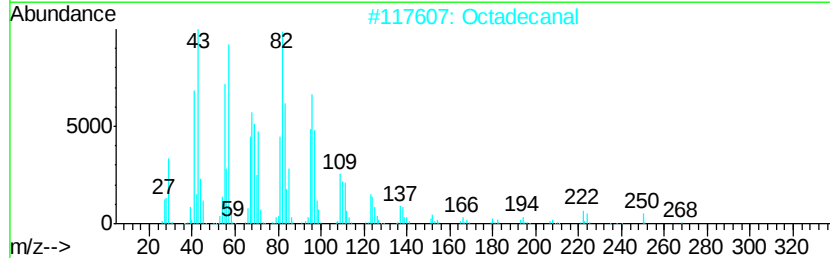
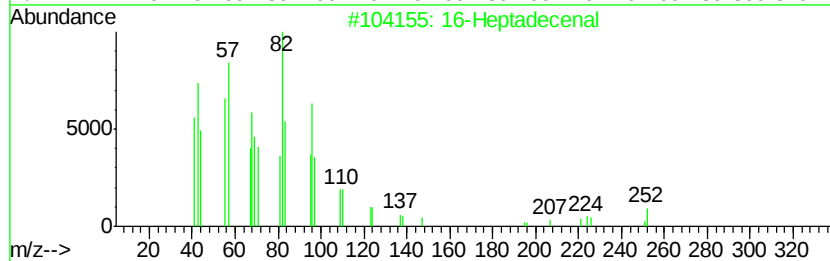
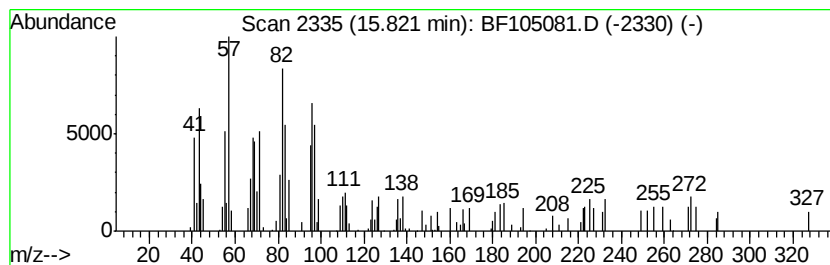
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 16-Heptadecenal Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	2.02 ng	67674	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			16-Heptadecenal	252	C17H32O	1000144-57-9	93
2			Octadecanal	268	C18H36O	000638-66-4	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			14-Heptadecenal	252	C17H32O	1000144-58-0	89
5			Hexadecanal	240	C16H32O	000629-80-1	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
Data File : BF105081.D
Acq On : 3 May 2018 19:10
Operator : JU/SJ
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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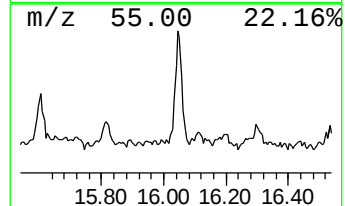
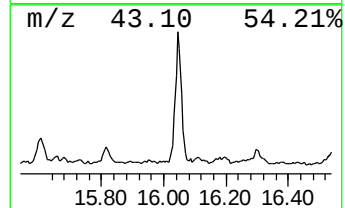
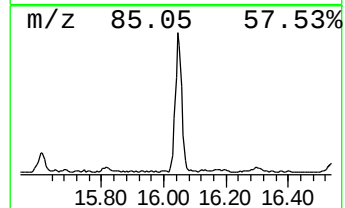
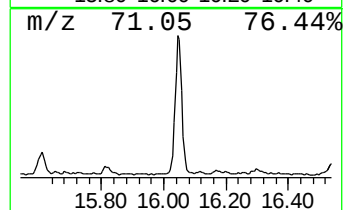
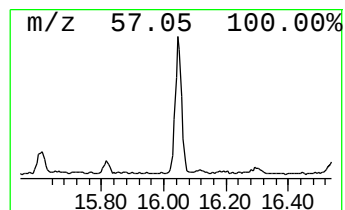
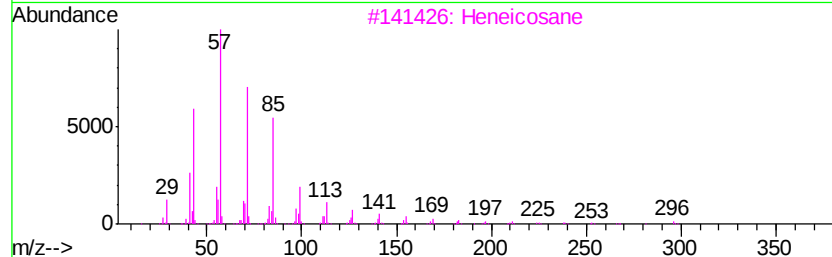
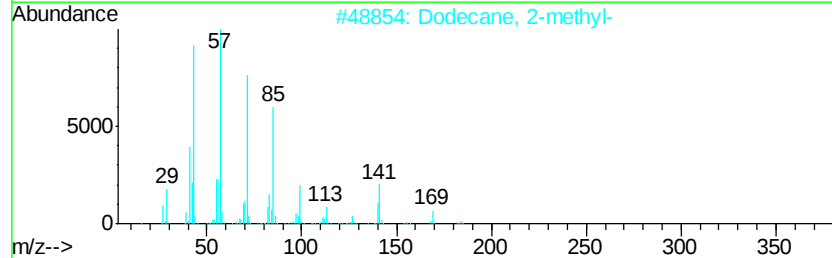
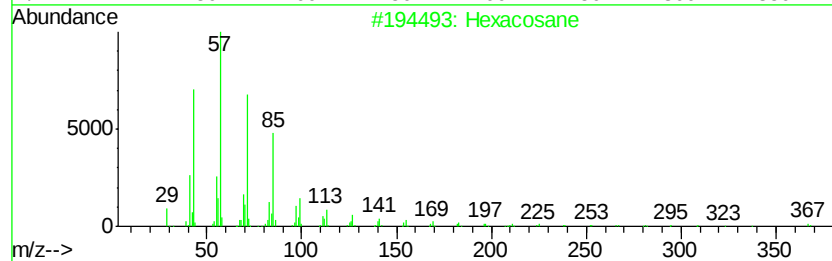
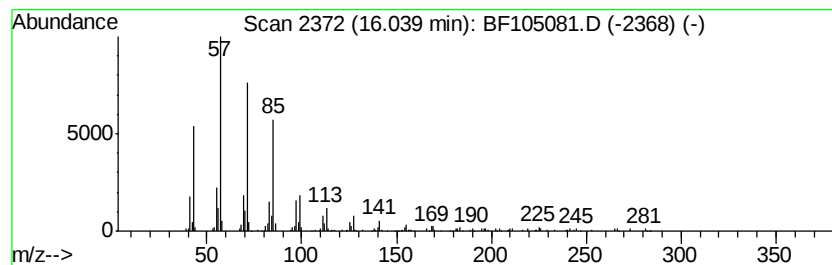
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	12.94 ng	432404	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		2-methyloctacosane	408	C29H60	1000376-72-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
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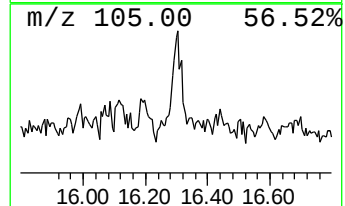
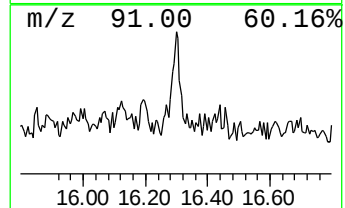
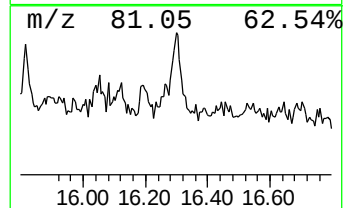
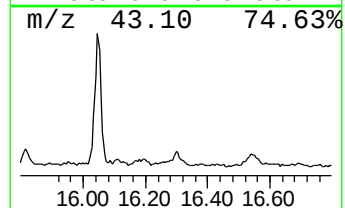
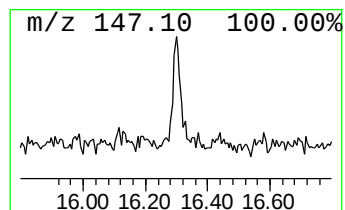
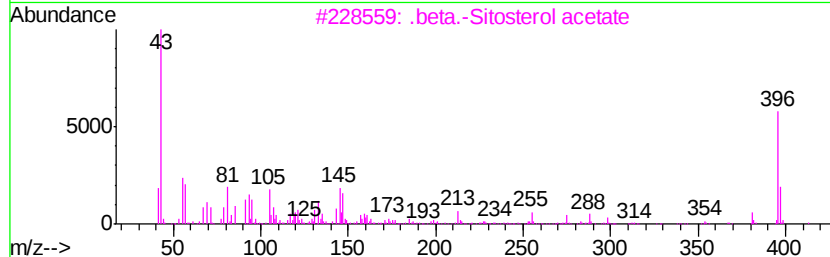
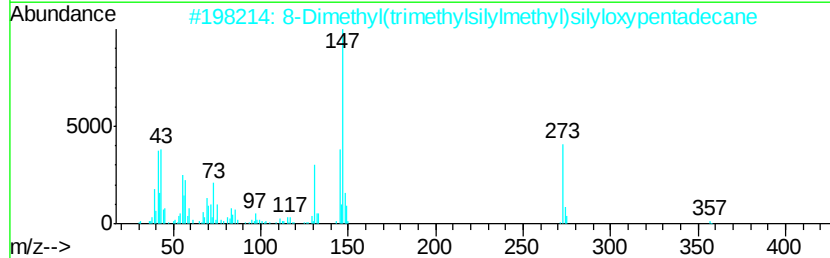
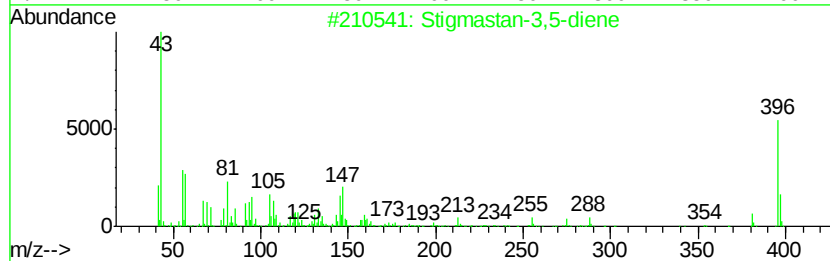
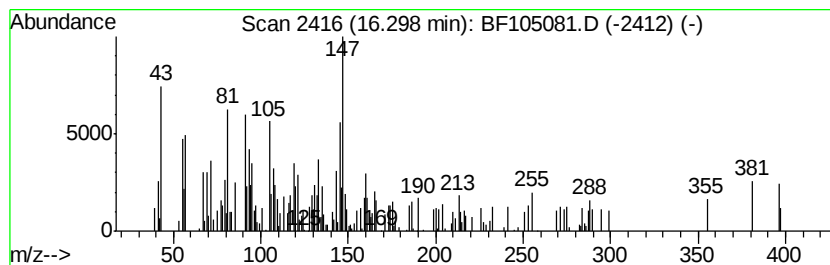
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Stigmastan-3,5-diene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.30	5.31 ng	177309	Perylene-d12	15.60		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Stigmastan-3,5-diene	396	C29H48	1000214-16-4	90
2		8-Dimethyl(trimethylsilylmethyl)...	372	C21H48OSi2	1000245-46-2	43
3		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	35
4		cis-1,1'-Bibenzocyclohexanylidene	288	C22H24	1000284-45-5	25
5		4-Methyl-1-pentamethyldisilyloxy...	232	C11H28OSi2	1000216-89-0	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
Data File : BF105081.D
Acq On : 3 May 2018 19:10
Operator : JU/SJ
Sample : J2665-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ES-RT-1871

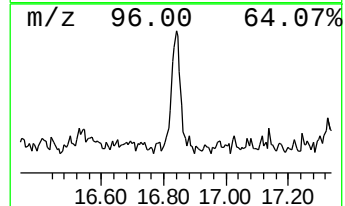
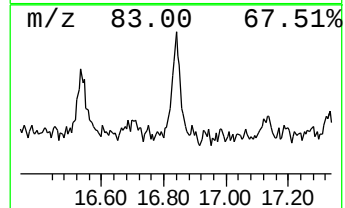
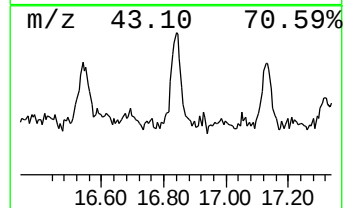
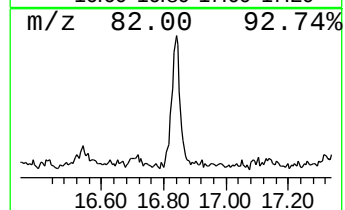
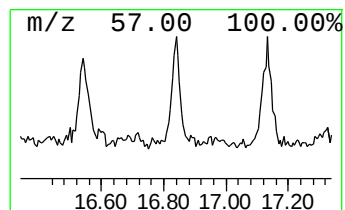
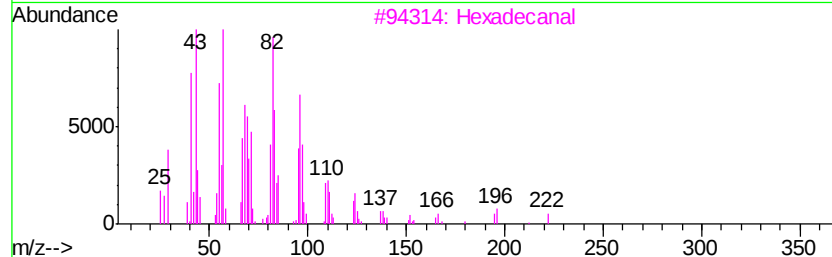
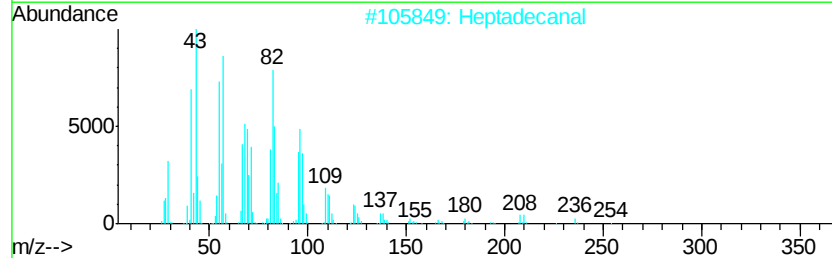
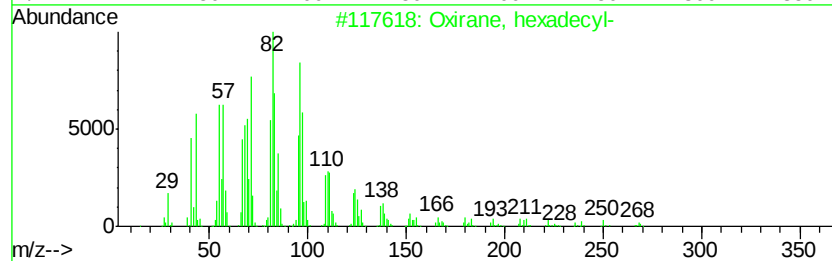
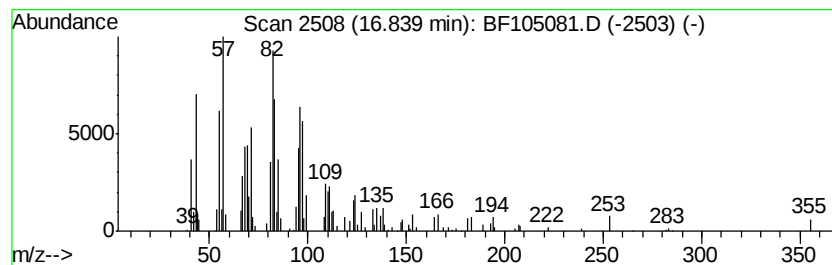
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Oxirane, hexadecyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.84	3.44 ng	114929	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2		Heptadecanal	254	C17H34O	1000376-70-0	90
3		Hexadecanal	240	C16H32O	000629-80-1	86
4		1,19-Eicosadiene	278	C20H38	014811-95-1	64
5		1,16-Hexadecanediol	258	C16H34O2	007735-42-4	64



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Sample : J2665-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ES-RT-1871

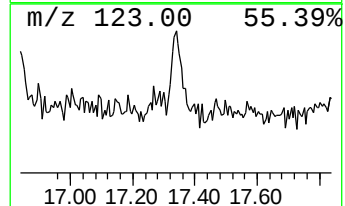
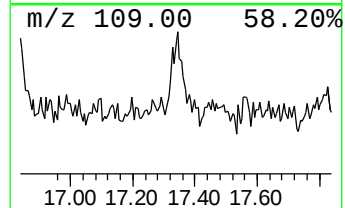
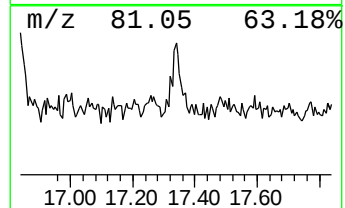
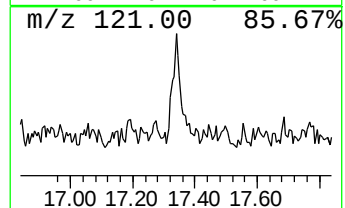
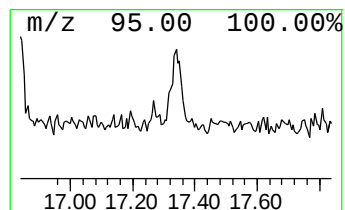
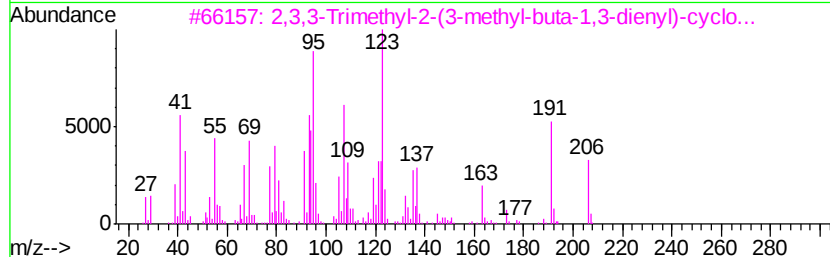
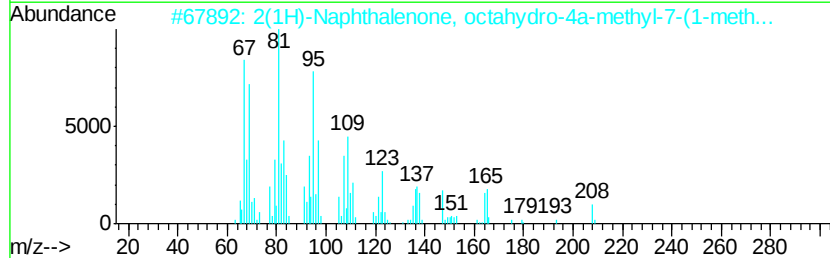
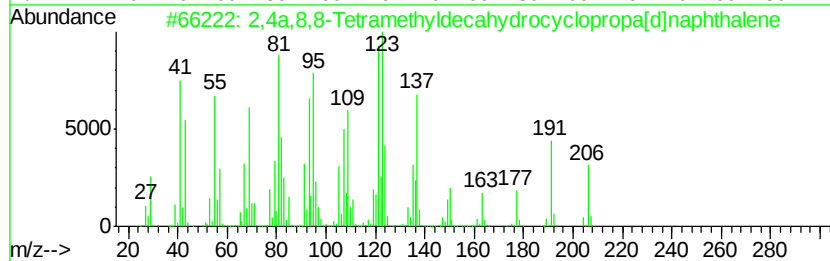
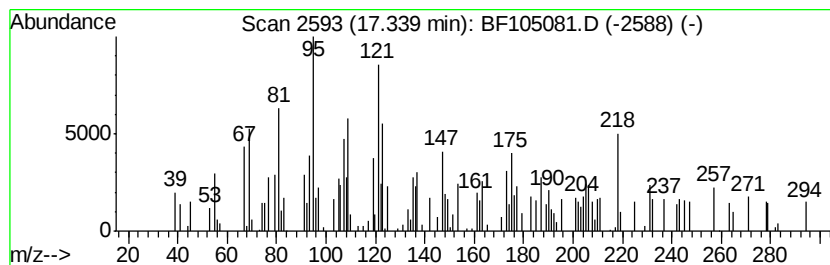
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.34 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	3.34 ng	111520	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	45
2		2(1H)-Naphthalenone, octahydro-4...	208	C14H24O	054594-42-2	38
3		2,3,3-Trimethyl-2-(3-methyl-buta...	206	C14H22O	1000193-60-6	38
4		2(1H)-Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	38
5		Acetic acid, trifluoro-, 3,4-dim...	218	C10H9F3O2	001957-55-7	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050318\
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 Acq On : 3 May 2018 19:10
 Operator : JU/SJ
 Sample : J2665-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ES-RT-1871

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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
2-Pentanone, 4-hy...	5.00	3.2 ng		92127	1	6.77	580597	20.0
unknown6.56	6.56	81.6 ng		2369700	1	6.77	580597	20.0
n-Hexadecanoic acid	11.84	10.4 ng		456458	4	11.31	875994	20.0
Octadecanoic acid	12.62	3.1 ng		135808	4	11.31	875994	20.0
Heptadecyl triflu...	13.83	6.6 ng		195764	5	14.02	593488	20.0
Erucic acid	14.12	2.7 ng		81222	5	14.02	593488	20.0
2-Octyldodecyl bu...	14.17	2.2 ng		66384	5	14.02	593488	20.0
Heneicosane	14.53	4.1 ng		121943	5	14.02	593488	20.0
1-Dodecanol, 2-oc...	14.87	3.4 ng		112590	6	15.60	668433	20.0
16-Heptadecenal	15.82	2.0 ng		67674	6	15.60	668433	20.0
Hexacosane	16.04	12.9 ng		432404	6	15.60	668433	20.0
Stigmastan-3,5-diene	16.30	5.3 ng		177309	6	15.60	668433	20.0
Oxirane, hexadecyl-	16.84	3.4 ng		114929	6	15.60	668433	20.0
unknown17.34	17.34	3.3 ng		111520	6	15.60	668433	20.0