

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
 Data File : BF118070.D
 Acq On : 3 Dec 2019 18:57
 Operator : JU
 Sample : K6090-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-20-S

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-BF111319.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	25	rVB	889368	1181797	21.45%	2.706%
2	5.093	507	511	524	rVB	552326	768074	13.94%	1.759%
3	5.498	574	580	583	rBV	3892635	3764479	68.34%	8.621%
4	6.510	746	752	755	rBV	3552123	4002991	72.67%	9.167%
5	6.645	770	775	779	rBV	4167942	4221259	76.63%	9.667%
6	6.875	811	814	818	rBV	1090095	832373	15.11%	1.906%
7	7.034	837	841	844	rBV	4014122	3331650	60.48%	7.630%
8	7.439	905	910	913	rBV	2610329	2543024	46.17%	5.824%
9	8.163	1029	1033	1041	rBV	1255699	1063012	19.30%	2.434%
10	9.245	1212	1217	1220	rBV	5508552	4910913	89.15%	11.247%
11	9.633	1279	1283	1287	rBV	702468	555464	10.08%	1.272%
12	9.928	1328	1333	1336	rBV	1340060	1249408	22.68%	2.861%
13	10.722	1463	1468	1471	rBV	3659511	3479617	63.17%	7.969%
14	11.416	1582	1586	1589	rBV	1474126	1320136	23.97%	3.023%
15	11.939	1667	1675	1688	rBV	451976	478410	8.69%	1.096%
16	12.633	1790	1793	1799	rBV	104553	104611	1.90%	0.240%
17	12.727	1806	1809	1818	rBV	139960	147219	2.67%	0.337%
18	12.863	1828	1832	1834	rBV	94976	103938	1.89%	0.238%
19	12.886	1834	1836	1846	rVB2	91975	124512	2.26%	0.285%
20	13.010	1852	1857	1861	rBV	5216170	5508299	100.00%	12.615%
21	13.580	1948	1954	1956	rBV2	69320	68828	1.25%	0.158%
22	13.904	2005	2009	2020	rVB2	289910	383544	6.96%	0.878%
23	14.068	2030	2037	2040	rBV	1275589	1152538	20.92%	2.639%
24	14.227	2059	2064	2069	rBV	125331	118391	2.15%	0.271%
25	14.533	2112	2116	2122	rBV2	164187	199753	3.63%	0.457%
26	14.845	2165	2169	2174	rBV	161581	168826	3.06%	0.387%
27	15.121	2211	2216	2220	rBV	36536	61395	1.11%	0.141%
28	15.192	2224	2228	2232	rBV	171310	197544	3.59%	0.452%
29	15.562	2285	2291	2305	rBV	784477	1239489	22.50%	2.839%
30	16.039	2366	2372	2377	rBV	108552	168185	3.05%	0.385%
31	16.557	2455	2460	2466	rBV3	56667	98931	1.80%	0.227%
32	17.427	2603	2608	2616	rVB8	30419	61198	1.11%	0.140%
33	17.680	2644	2651	2656	rBV10	21738	55844	1.01%	0.128%

LSC Area Percent Report

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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-20-S

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-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

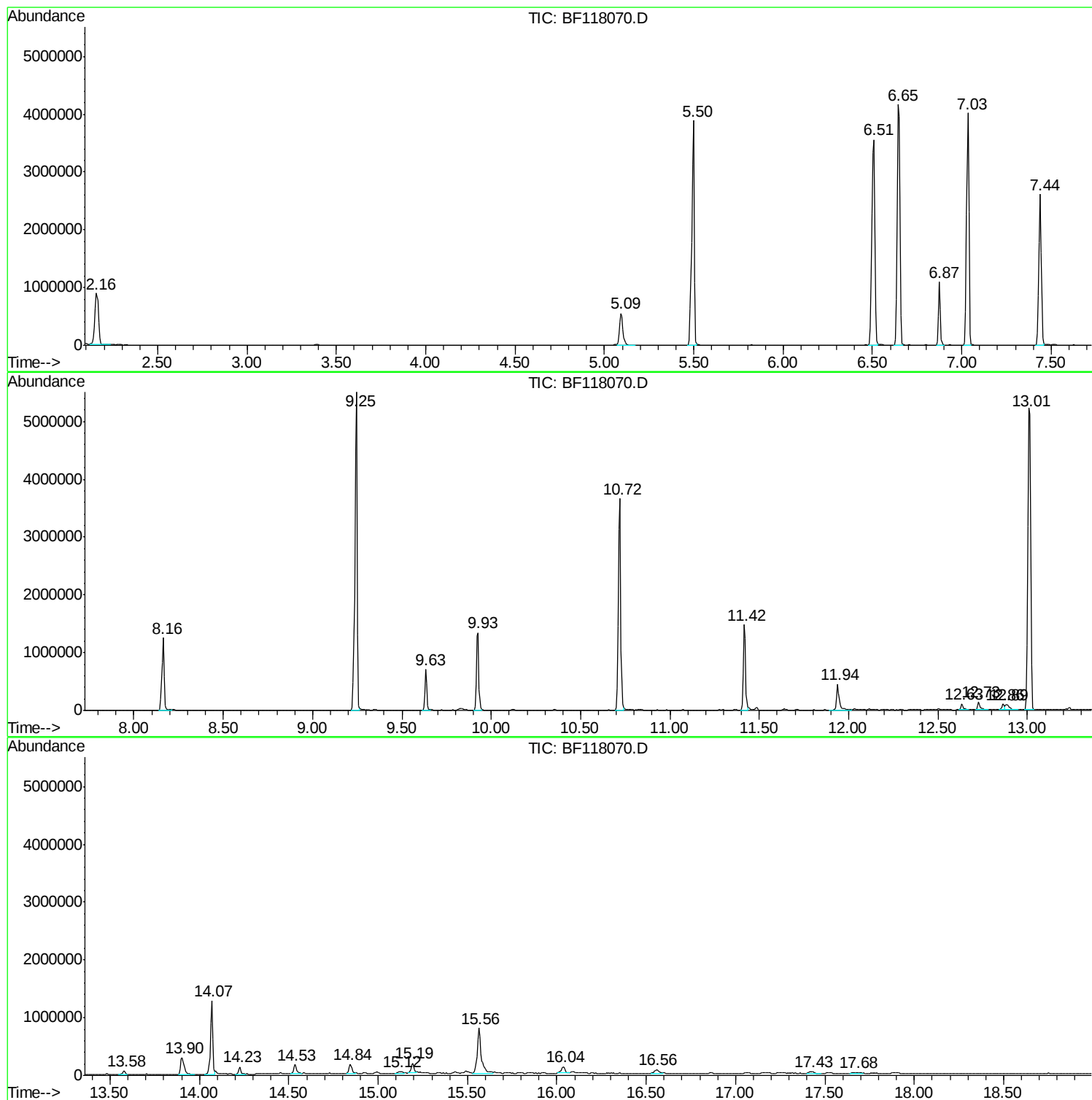
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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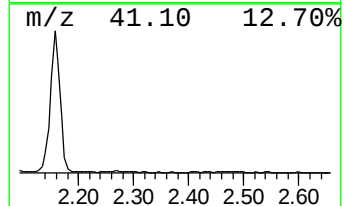
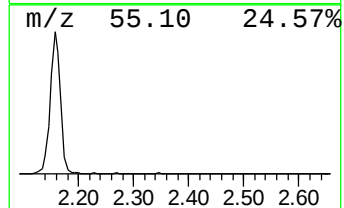
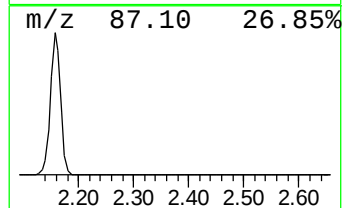
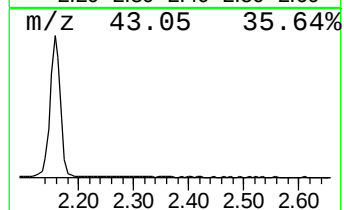
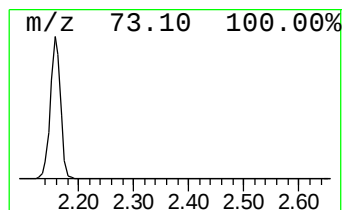
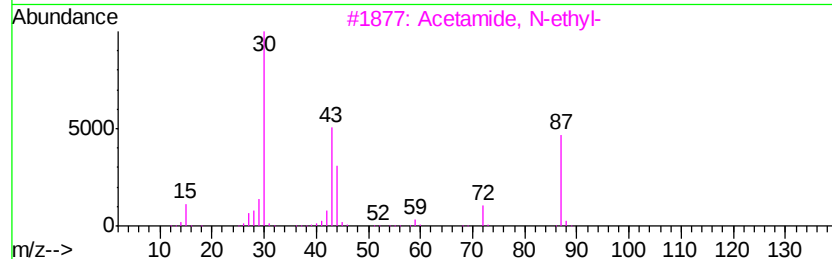
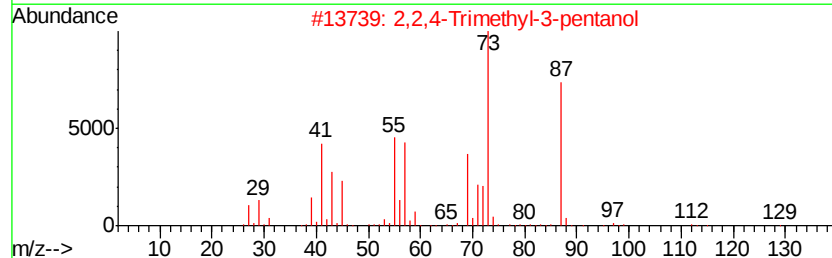
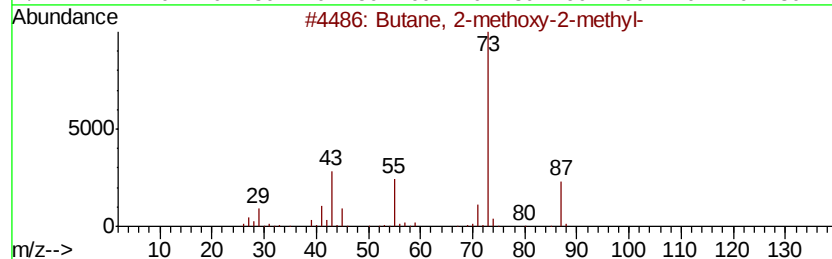
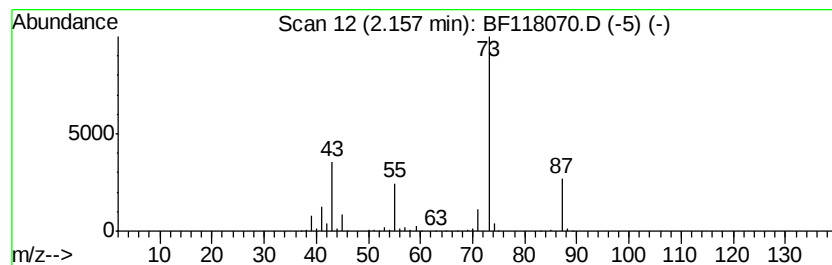
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	28.40 ng	1181800	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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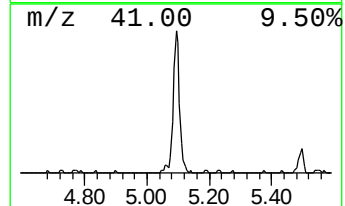
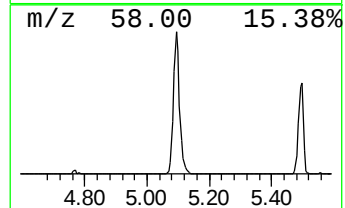
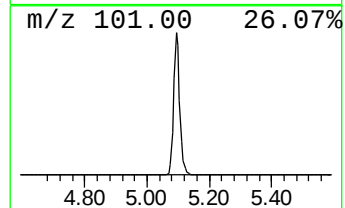
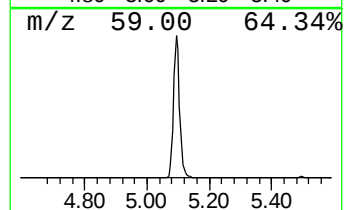
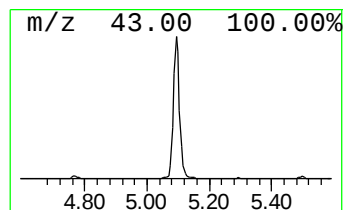
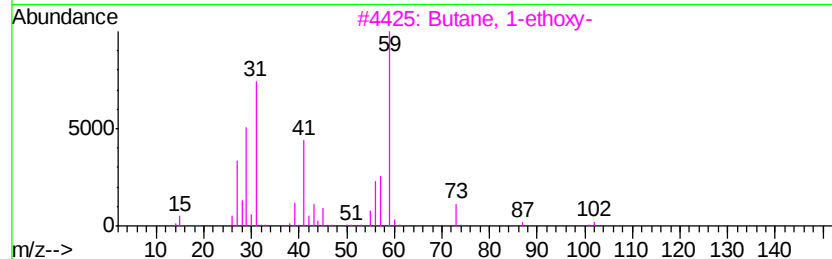
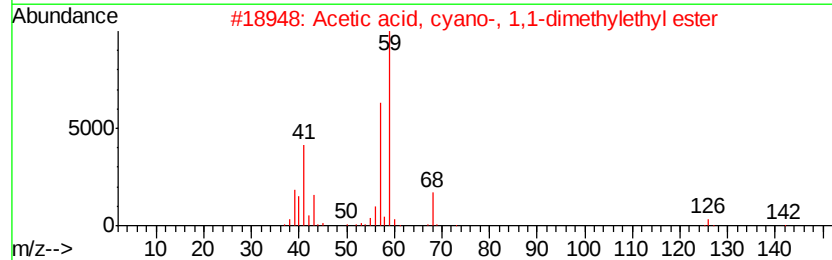
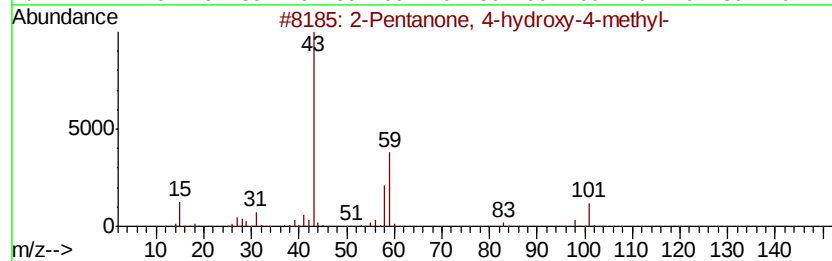
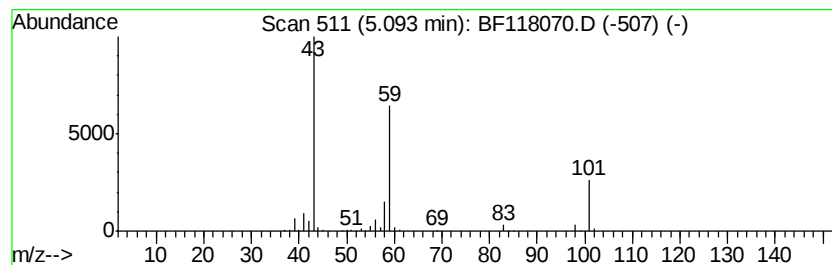
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	18.46 ng	768074	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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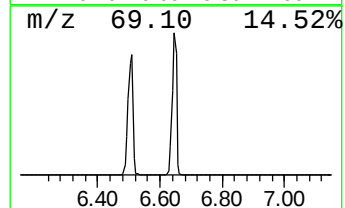
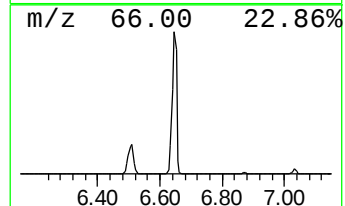
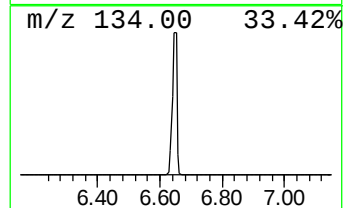
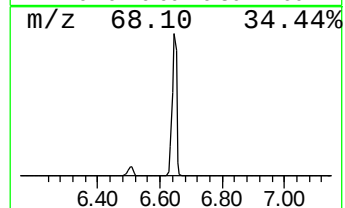
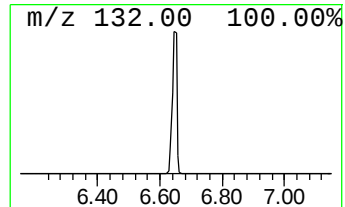
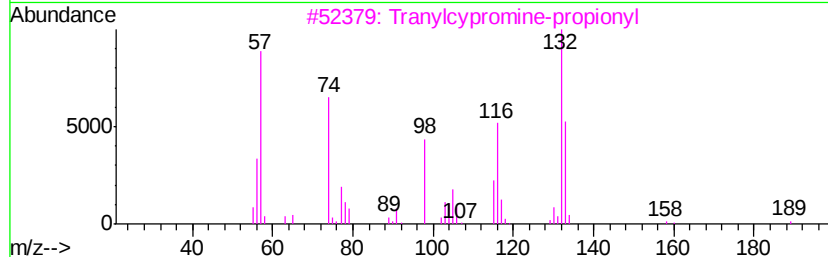
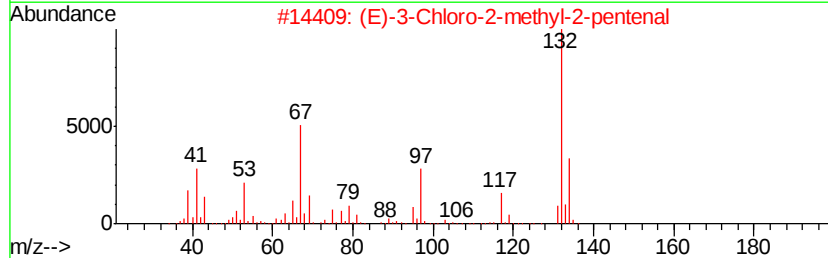
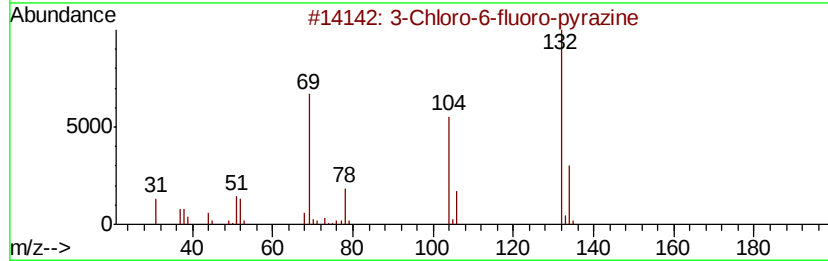
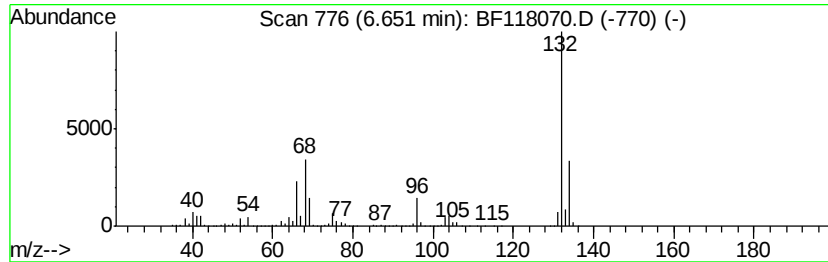
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	101.43 ng	4221260	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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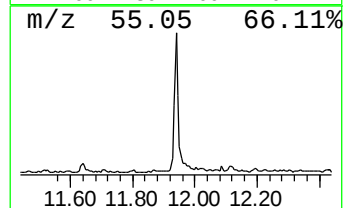
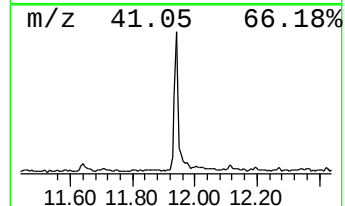
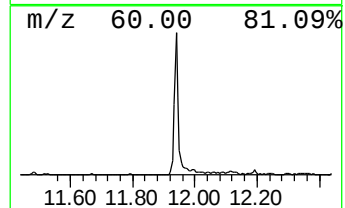
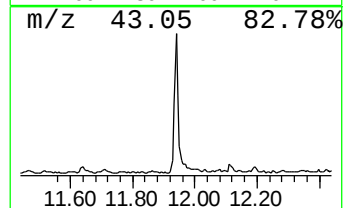
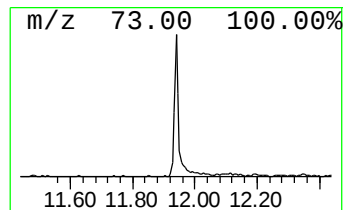
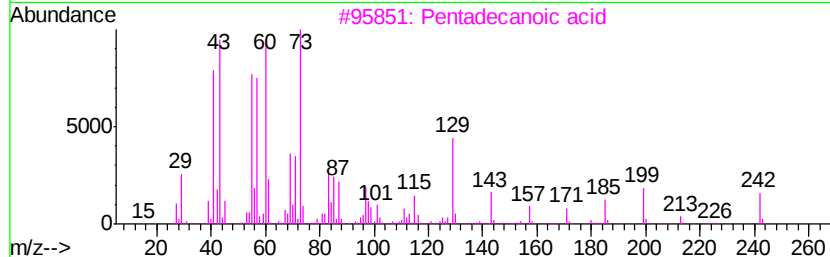
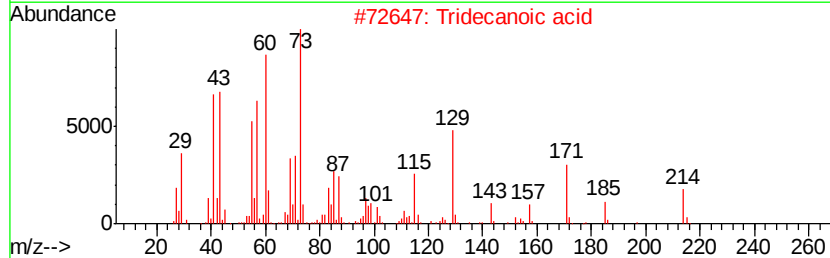
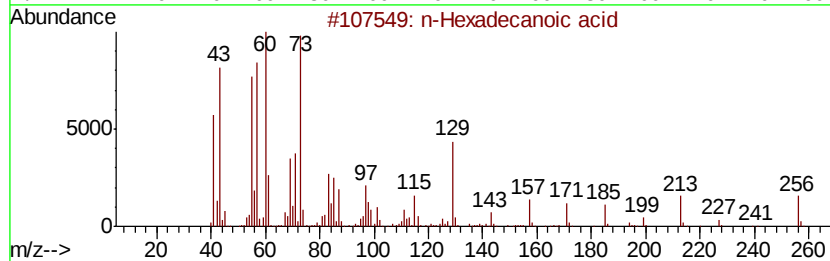
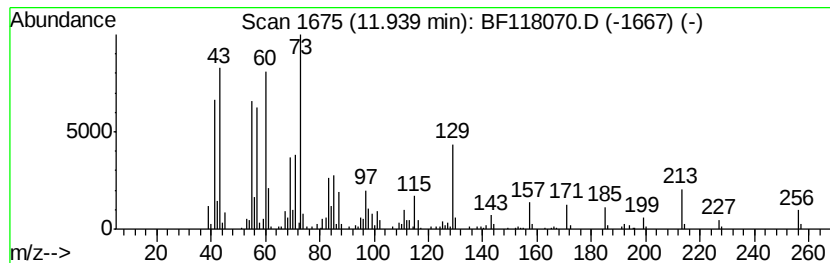
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	7.25 ng	478410	Phenanthrene-d10	11.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	60



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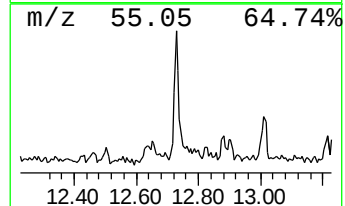
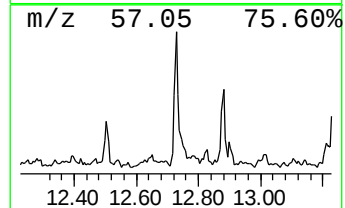
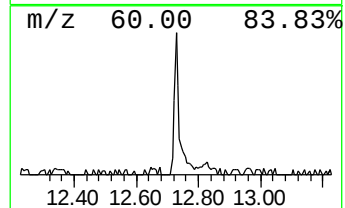
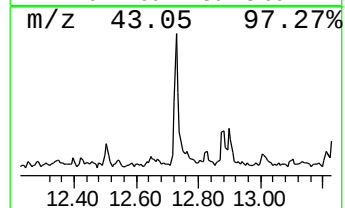
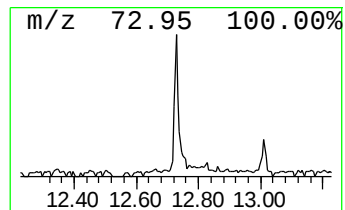
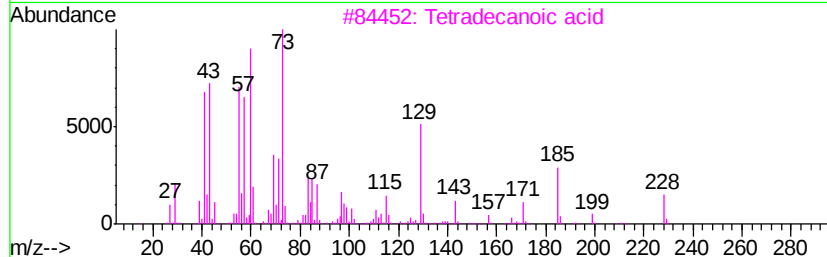
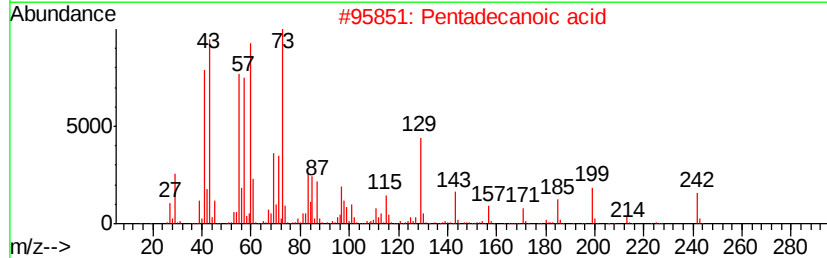
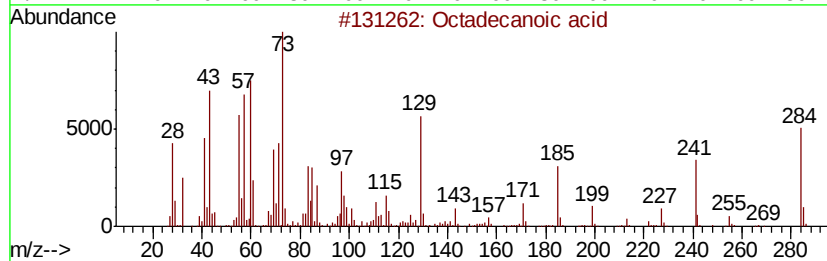
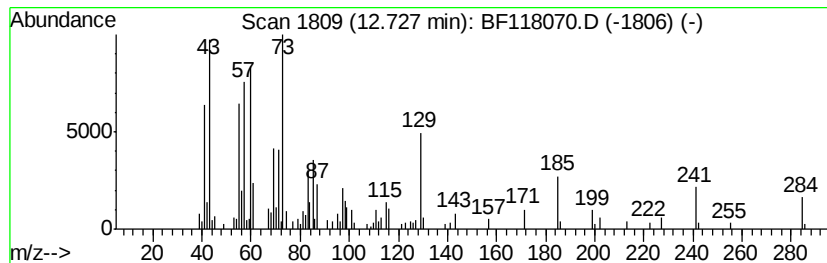
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.73	2.23 ng	147219	Phenanthrene-d10		11.42
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	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5	Undecanoic acid	186	C11H22O2	000112-37-8	43



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Misc :
ALS Vial : 13 Sample Multiplier: 1

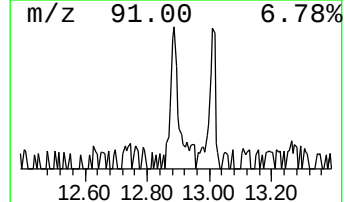
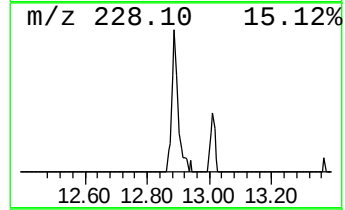
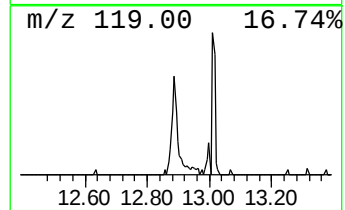
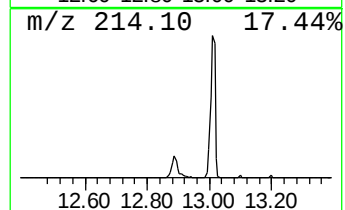
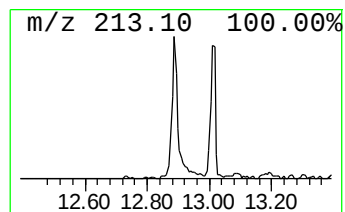
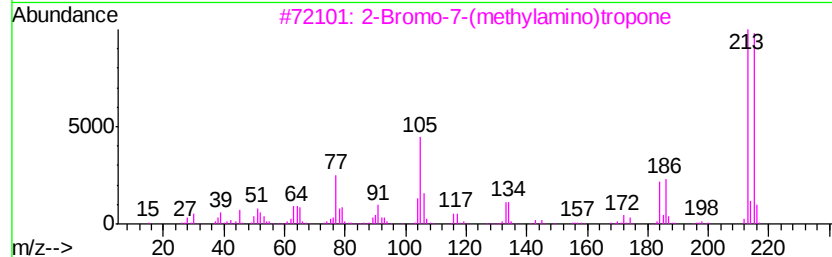
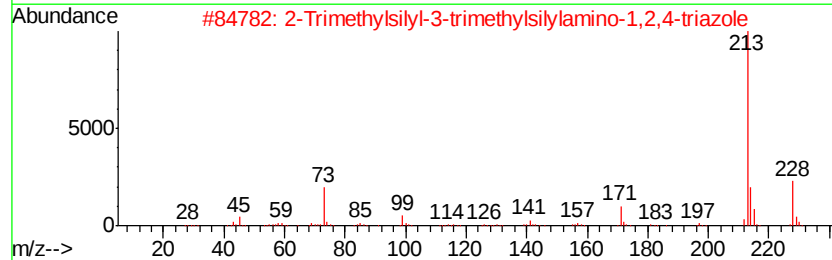
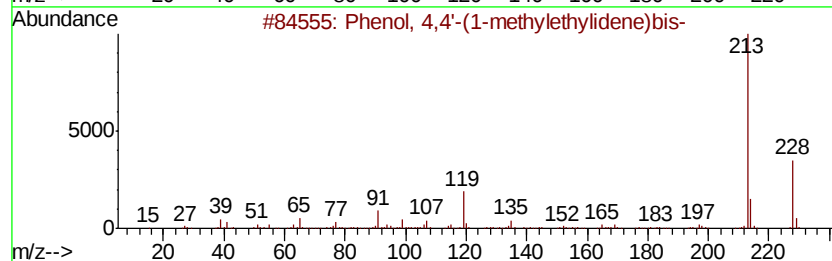
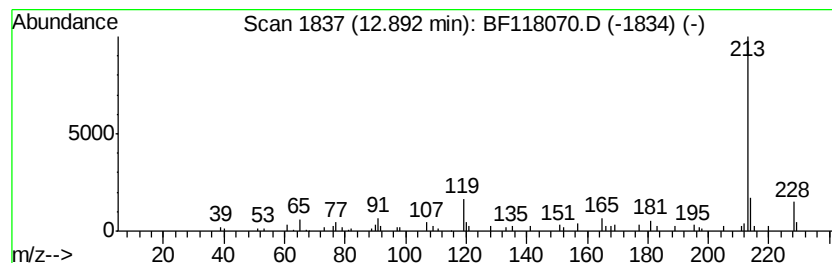
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ClientSampleId :
TP-20-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.89	2.16 ng	124512	Chrysene-d12	14.07		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene)bis-	228	C15H16O2	000080-05-7	91	
2	2-Trimethylsilyl-3-trimethylsilyl-	228	C8H20N4Si2	1000373-05-5	64	
3	2-Bromo-7-(methylamino)tropon	213	C8H8BrNO	103539-25-9	50	
4	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	50	
5	2-Trifluoromethyl-5-hydroxyquino...	213	C10H6F3NO	041958-06-9	50	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118070.D
Acq On : 3 Dec 2019 18:57
Operator : JU
Sample : K6090-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-20-S

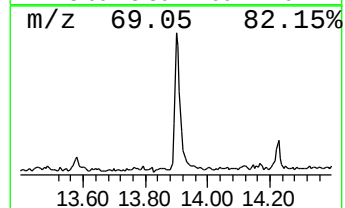
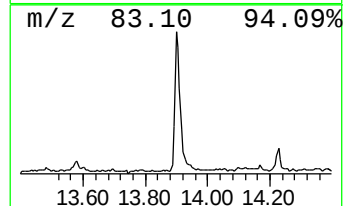
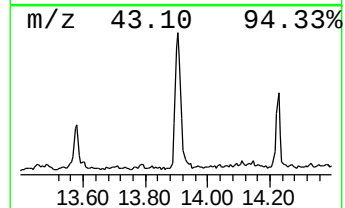
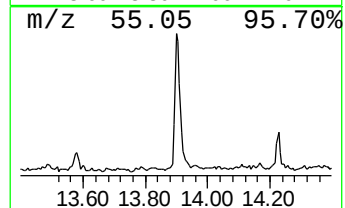
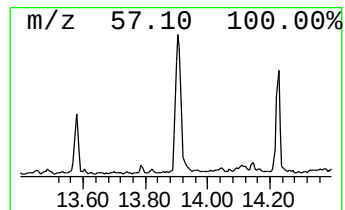
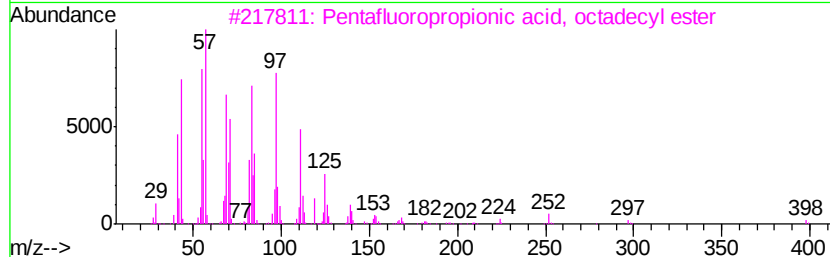
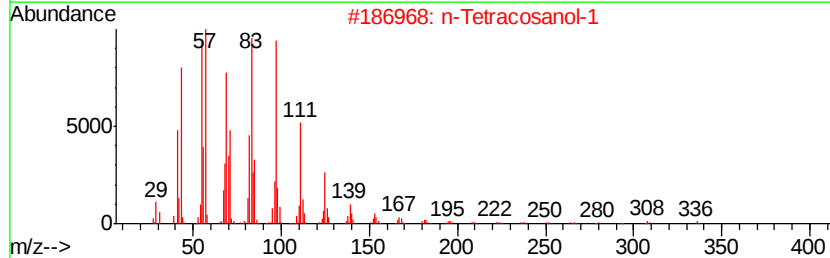
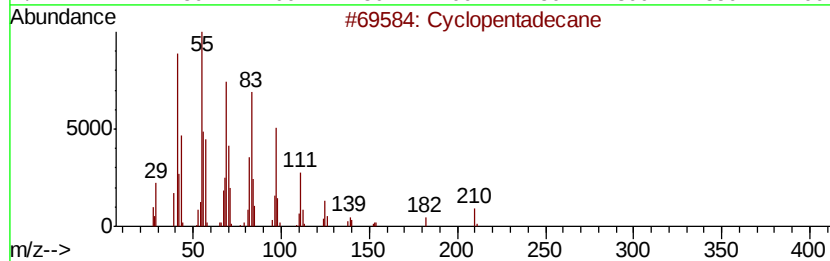
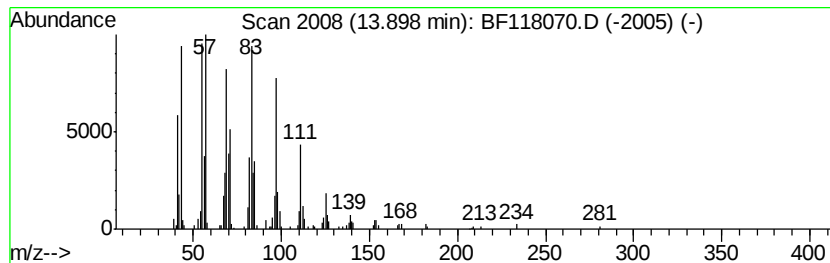
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	6.66 ng	383544	Chrysene-d12	14.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentadecane	210	C15H30	000295-48-7	96
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
3		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	91
4		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	91
5		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
 Data File : BF118070.D
 Acq On : 3 Dec 2019 18:57
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Instrument :
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 ClientSampleId :
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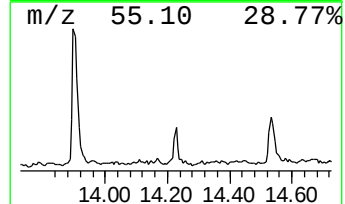
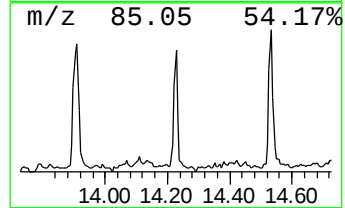
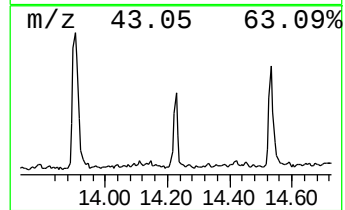
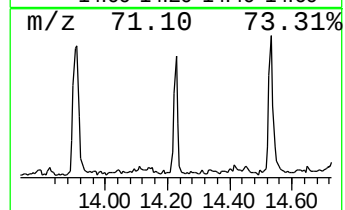
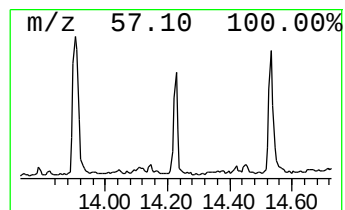
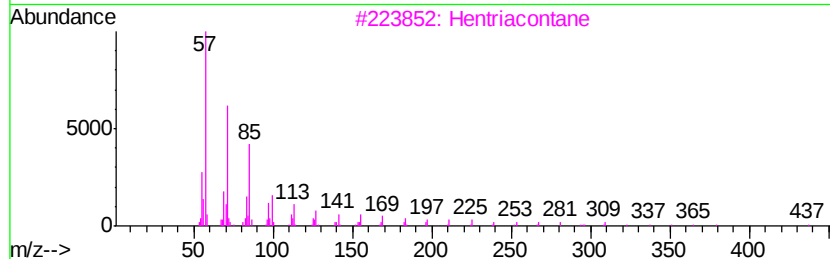
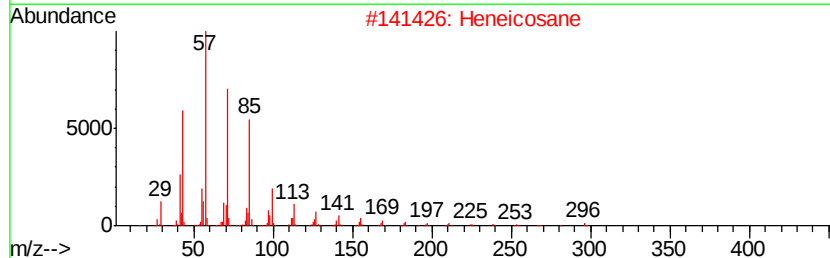
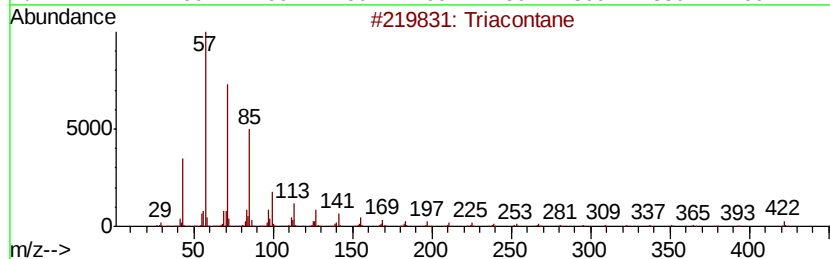
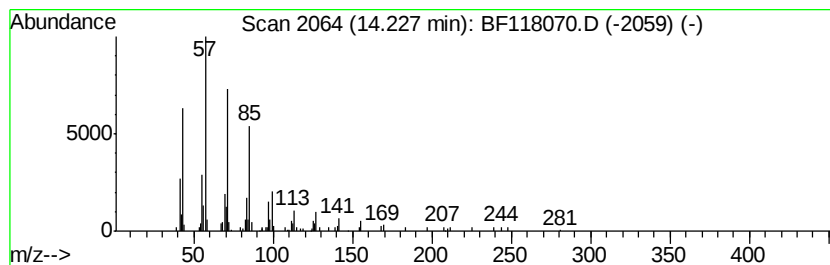
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Triacontane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	2.05 ng	118391	Chrysene-d12	14.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontane	422	C30H62	000638-68-6	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hentriacontane	437	C31H64	000630-04-6	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118070.D
Acq On : 3 Dec 2019 18:57
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Sample : K6090-03
Misc :
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Instrument :
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ClientSampleId :
TP-20-S

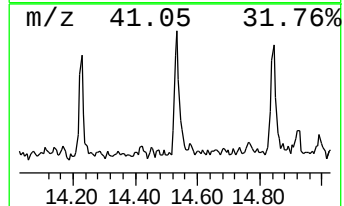
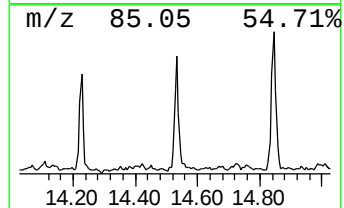
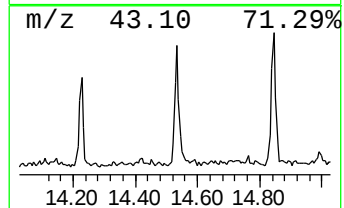
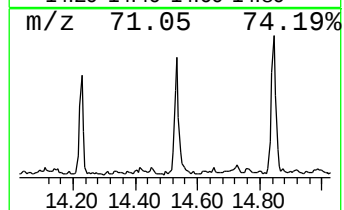
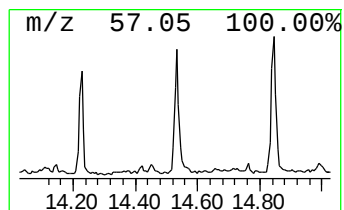
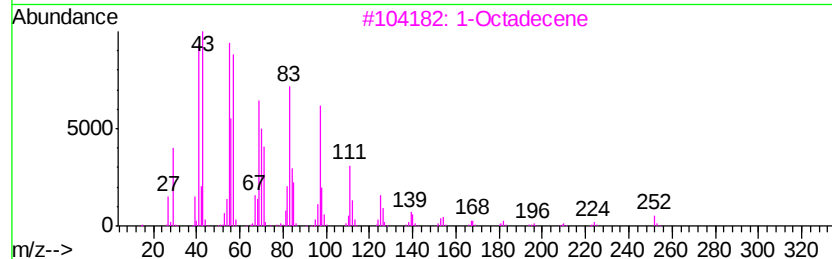
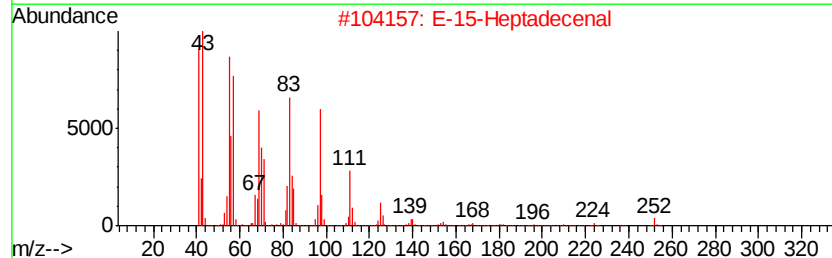
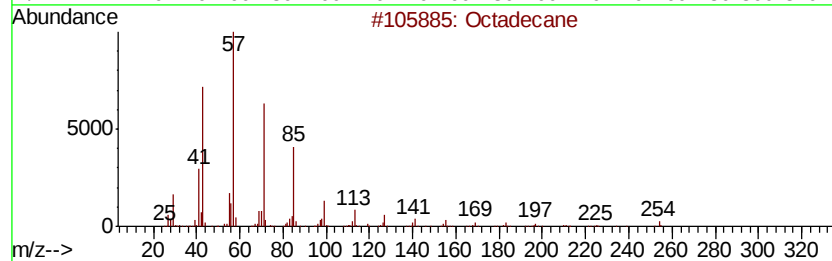
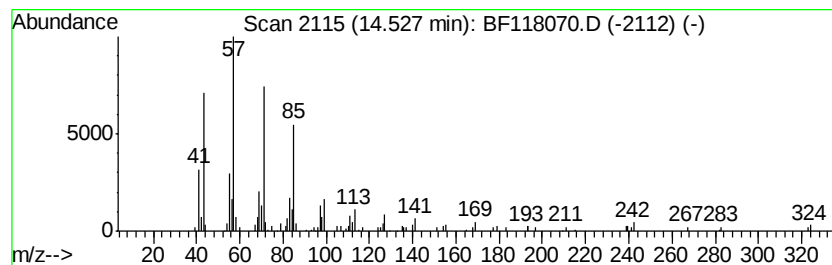
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	3.47 ng	199753	Chrysene-d12	14.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	98
2			E-15-Heptadecenal	252	C17H32O	1000130-97-9	92
3			1-Octadecene	252	C18H36	000112-88-9	92
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Tritetracontane	605	C43H88	007098-21-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
Data File : BF118070.D
Acq On : 3 Dec 2019 18:57
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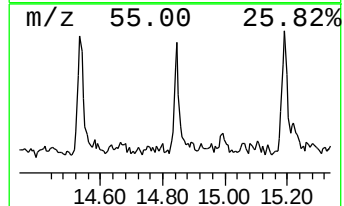
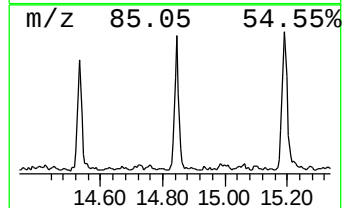
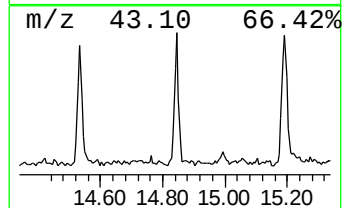
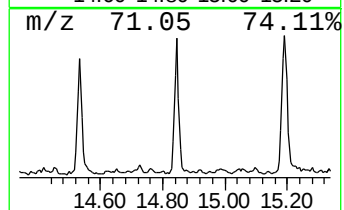
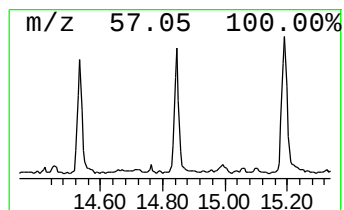
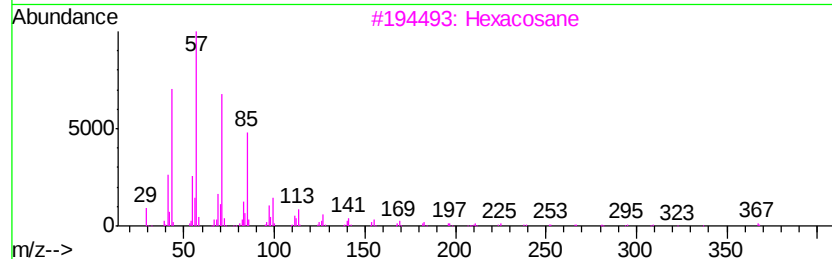
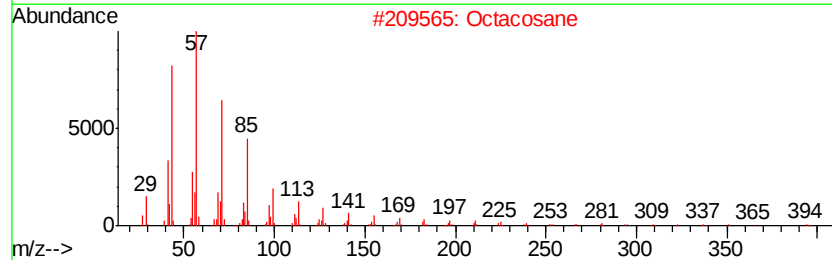
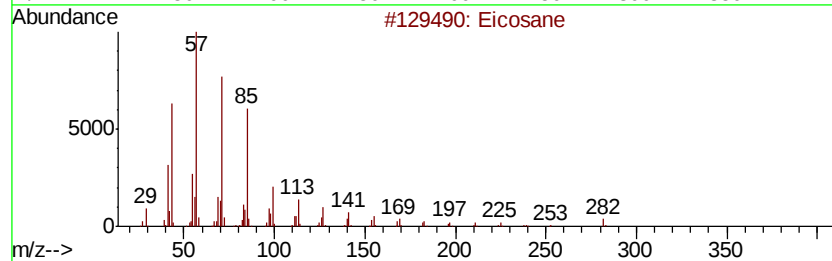
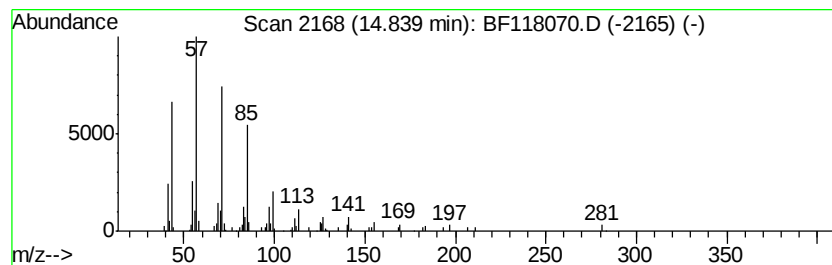
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	2.72 ng	168826	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
 Data File : BF118070.D
 Acq On : 3 Dec 2019 18:57
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Instrument :
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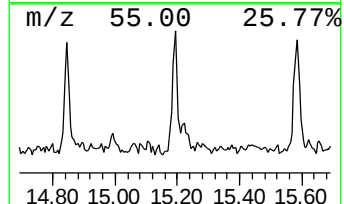
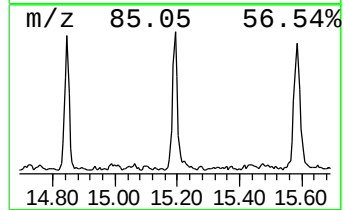
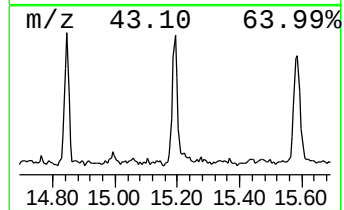
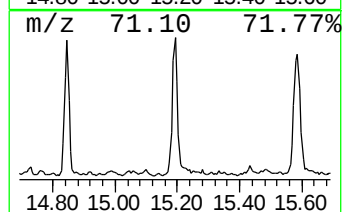
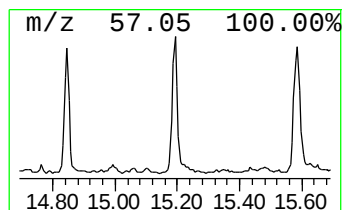
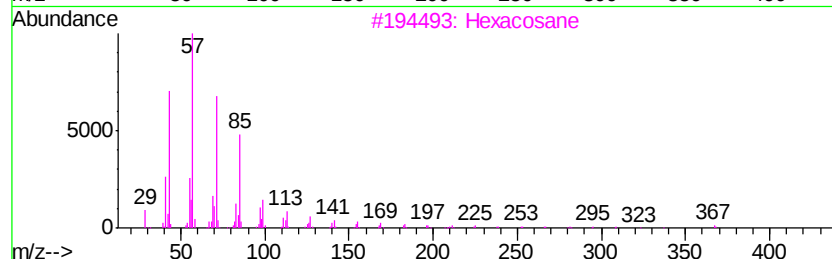
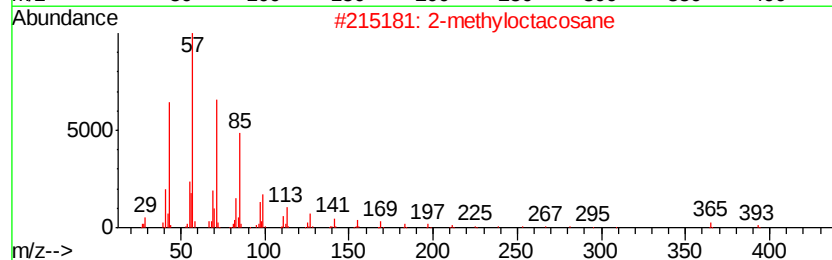
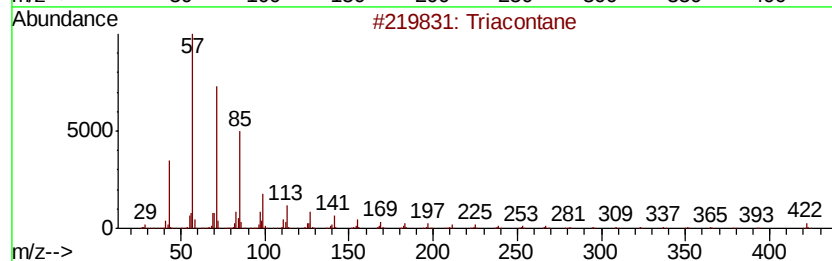
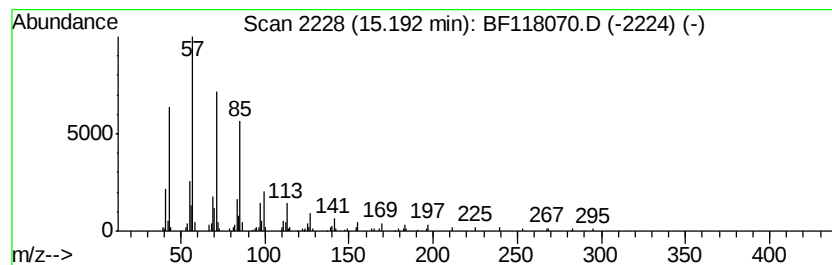
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 2-methyloctacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	3.19 ng	197544	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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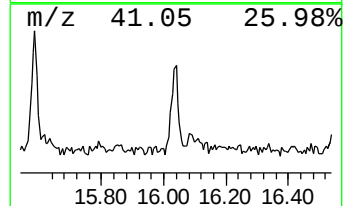
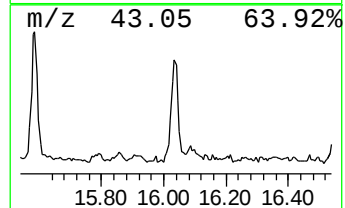
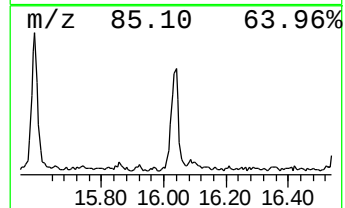
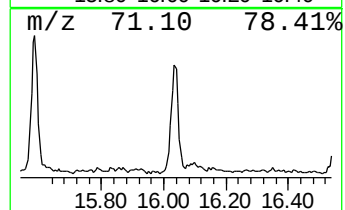
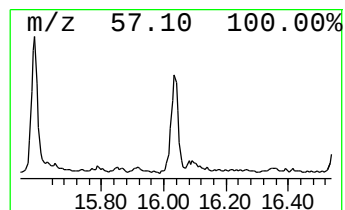
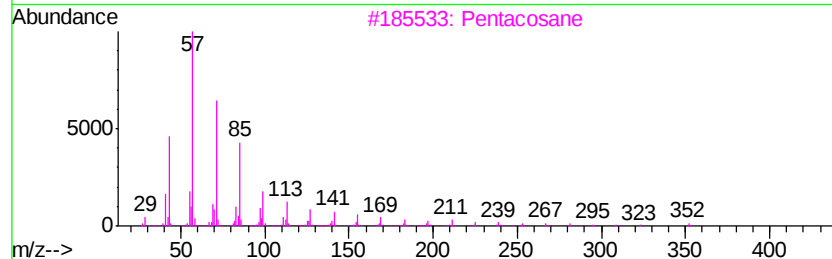
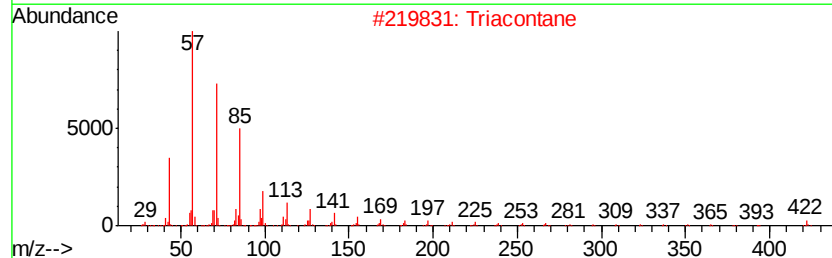
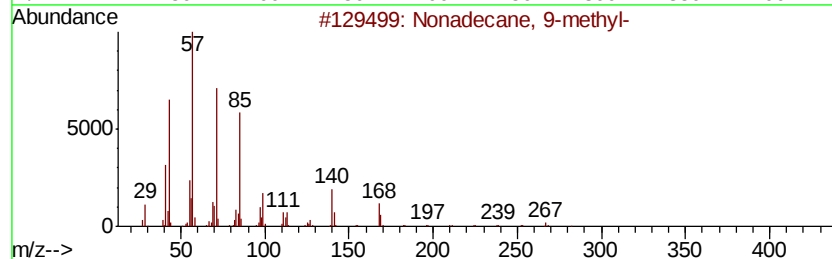
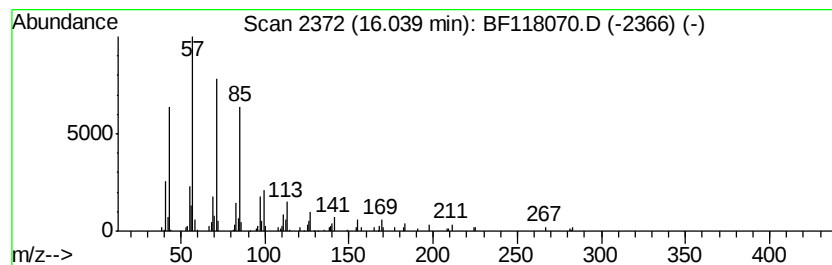
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonadecane, 9-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.71 ng	168185	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptadecane, 2,3-dimethyl-	268	C19H40	061868-03-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF120319\
 Data File : BF118070.D
 Acq On : 3 Dec 2019 18:57
 Operator : JU
 Sample : K6090-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-20-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.16	28.4	ng	1181800	1	6.87	832373	20.0
2-Pentanone, 4-hy...	5.09	18.5	ng	768074	1	6.87	832373	20.0
unknown6.65	6.65	101.4	ng	4221260	1	6.87	832373	20.0
n-Hexadecanoic acid	11.94	7.3	ng	478410	4	11.42	1320140	20.0
Octadecanoic acid	12.73	2.2	ng	147219	4	11.42	1320140	20.0
Phenol, 4,4'-(1-m...	12.89	2.2	ng	124512	5	14.07	1152540	20.0
Cyclopentadecane	13.90	6.7	ng	383544	5	14.07	1152540	20.0
Triacontane	14.23	2.0	ng	118391	5	14.07	1152540	20.0
Octadecane	14.53	3.5	ng	199753	5	14.07	1152540	20.0
Eicosane	14.84	2.7	ng	168826	6	15.56	1239490	20.0
2-methyloctacosane	15.19	3.2	ng	197544	6	15.56	1239490	20.0
Nonadecane, 9-met...	16.04	2.7	ng	168185	6	15.56	1239490	20.0