

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
 Data File : BF110802.D
 Acq On : 15 Nov 2018 21:35
 Operator : JU/SJ
 Sample : J5950-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSI-20-TW-1

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-BF110818.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	2.240	21	26	37	rVB	93030	150990	8.21%	0.962%
2	5.181	522	526	542	rBV	47286	76616	4.17%	0.488%
3	5.428	565	568	575	rBV	36914	37197	2.02%	0.237%
4	5.534	583	586	587	rBV	27571	22064	1.20%	0.141%
5	5.563	587	591	612	rVB	1756428	1759946	95.68%	11.208%
6	6.569	757	762	778	rBV	1782738	1827734	99.37%	11.640%
7	6.710	782	786	789	rBV	2145241	1839314	100.00%	11.714%
8	6.939	821	825	832	rBV	405501	350103	19.03%	2.230%
9	7.092	847	851	863	rVB	1465750	1409759	76.65%	8.978%
10	7.463	911	914	917	rBV	50748	41485	2.26%	0.264%
11	7.504	917	921	931	rBV	1124986	1132457	61.57%	7.212%
12	7.951	994	997	1005	rVB	41328	44942	2.44%	0.286%
13	8.222	1040	1043	1046	rBV	467563	412817	22.44%	2.629%
14	9.298	1222	1226	1236	rBV	1998833	1835607	99.80%	11.690%
15	9.686	1289	1292	1300	rBV	223640	215739	11.73%	1.374%
16	9.980	1338	1342	1353	rVB2	472900	441903	24.03%	2.814%
17	10.145	1367	1370	1375	rBV	33246	39655	2.16%	0.253%
18	10.369	1404	1408	1410	rBV	50633	54447	2.96%	0.347%
19	10.386	1410	1411	1419	rVB2	43980	42331	2.30%	0.270%
20	10.486	1425	1428	1439	rVB	72374	69712	3.79%	0.444%
21	10.774	1474	1477	1486	rBV	1026703	993520	54.02%	6.327%
22	10.863	1490	1492	1496	rBV	28998	31052	1.69%	0.198%
23	10.898	1496	1498	1503	rVV	40500	49891	2.71%	0.318%
24	10.974	1507	1511	1514	rVB3	27313	30325	1.65%	0.193%
25	11.474	1593	1596	1604	rBV	364627	356230	19.37%	2.269%
26	11.980	1679	1682	1687	rBV	59611	56710	3.08%	0.361%
27	13.063	1861	1866	1869	rBV	1418429	1293613	70.33%	8.238%
28	13.933	2011	2014	2024	rBV	53293	68460	3.72%	0.436%
29	14.127	2044	2047	2053	rBV	347019	365351	19.86%	2.327%
30	14.251	2064	2068	2071	rBV	23694	24230	1.32%	0.154%
31	14.557	2117	2120	2128	rBV	38733	54781	2.98%	0.349%
32	14.874	2170	2174	2178	rBV	45812	47596	2.59%	0.303%
33	15.221	2230	2233	2238	rBV	48768	54193	2.95%	0.345%
34	15.615	2296	2300	2303	rBV	44988	59506	3.24%	0.379%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.662	2303	2308	2317	rVB	186771	272228	14.80%	1.734%
36	16.074	2373	2378	2384	rVB2	40538	64548	3.51%	0.411%
37	16.604	2463	2468	2474	rVB3	27269	45168	2.46%	0.288%
38	17.221	2568	2573	2578	rBV5	13371	29944	1.63%	0.191%

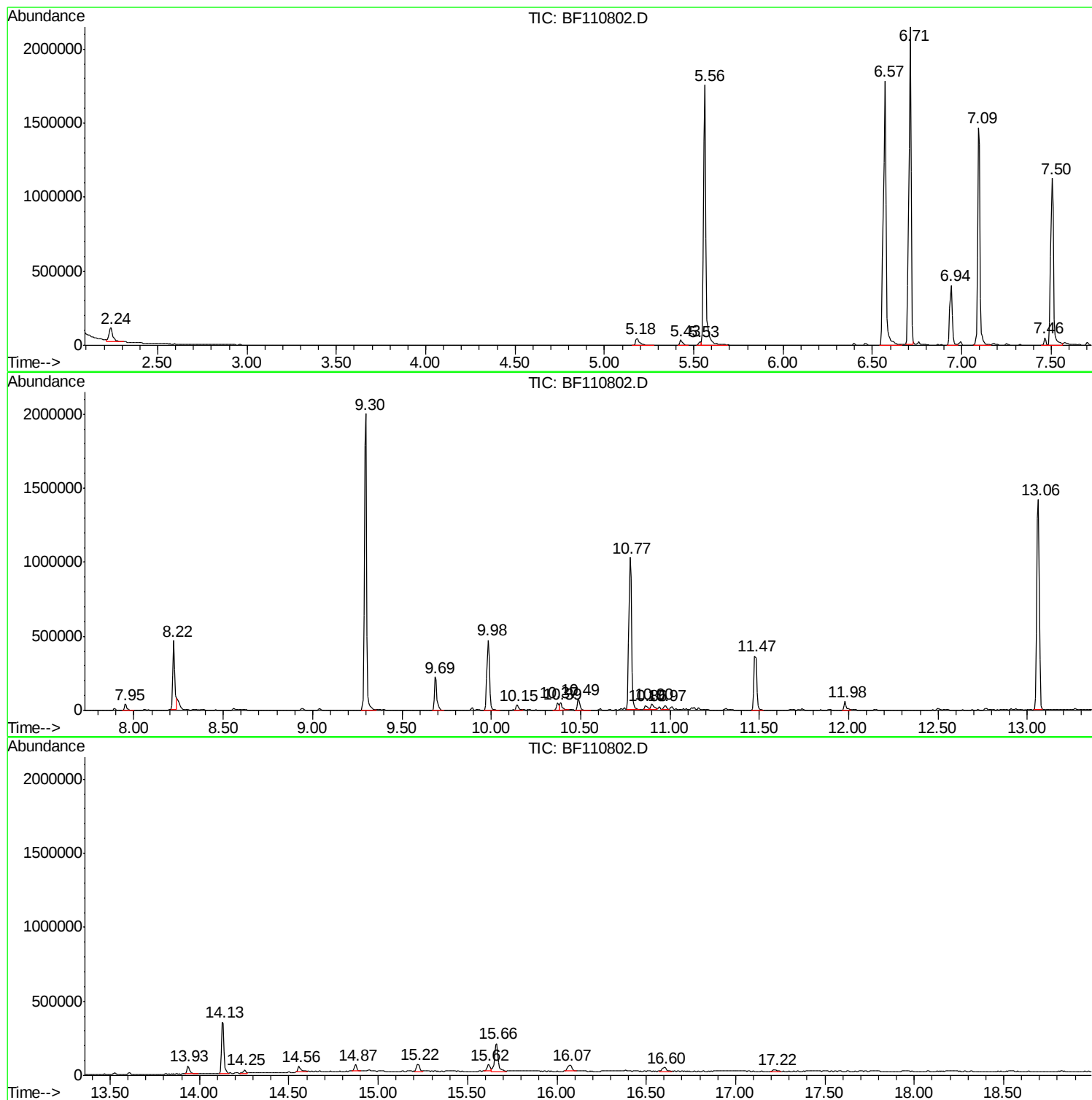
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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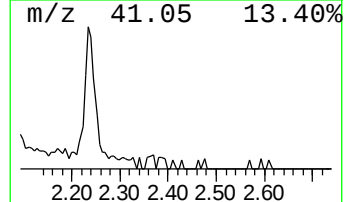
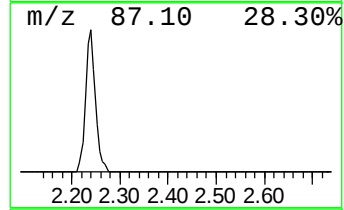
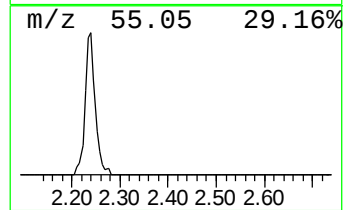
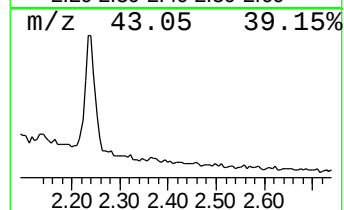
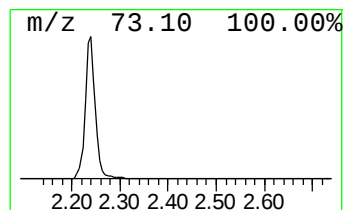
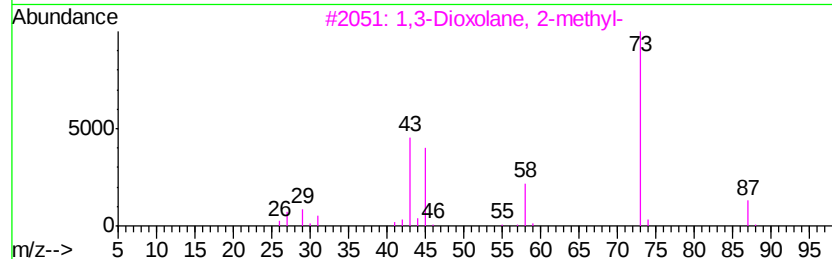
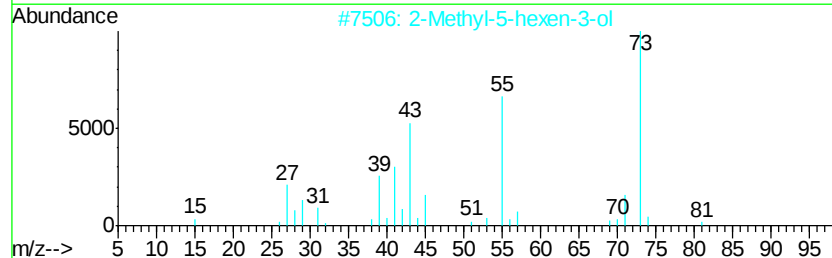
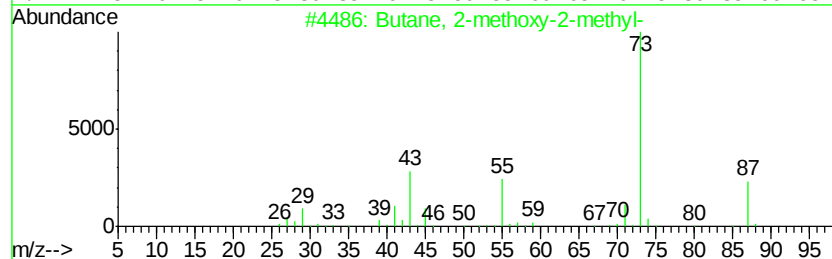
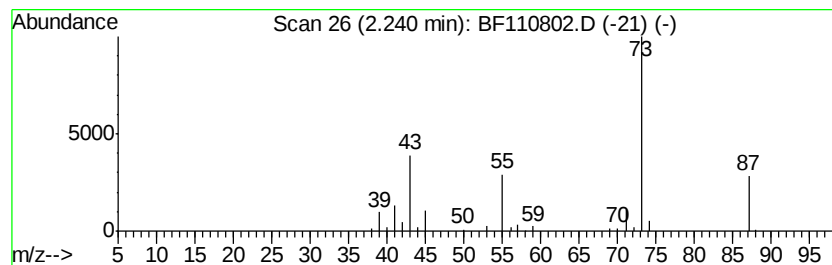
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.24	8.63 ng	150990	1,4-Dichlorobenzene-d4	6.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	38
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
5			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	25



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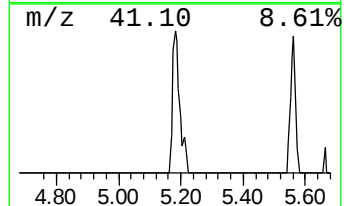
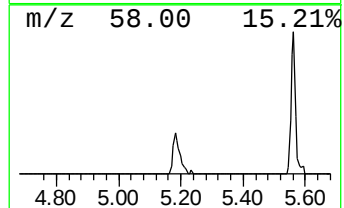
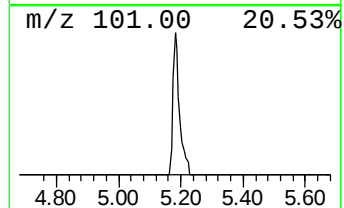
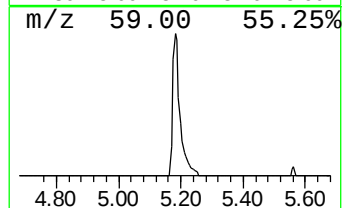
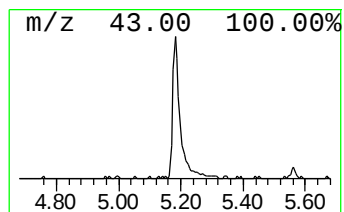
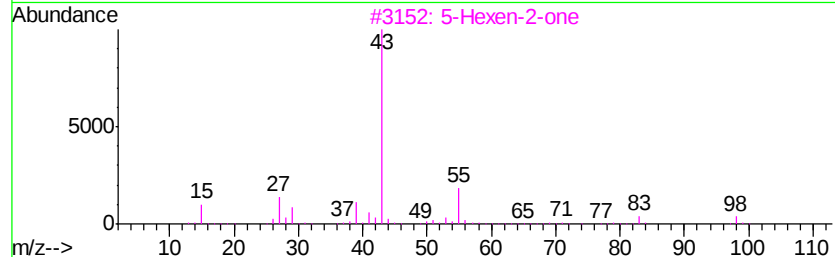
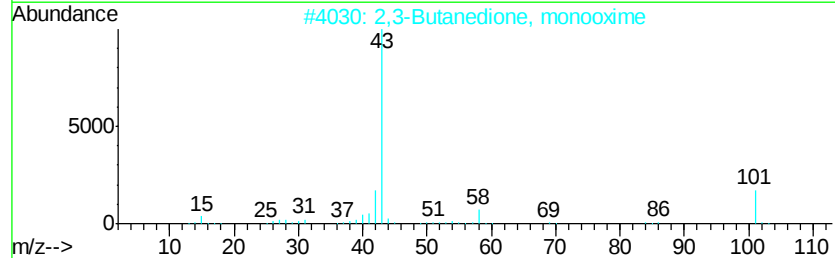
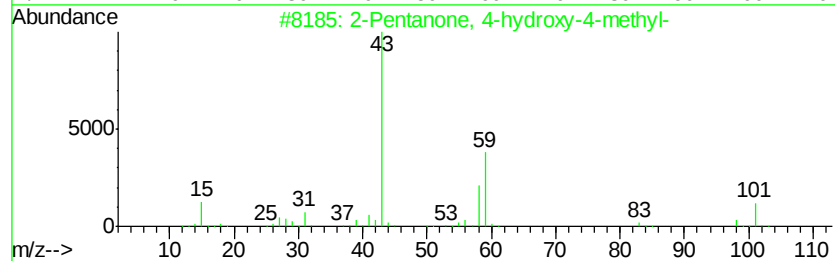
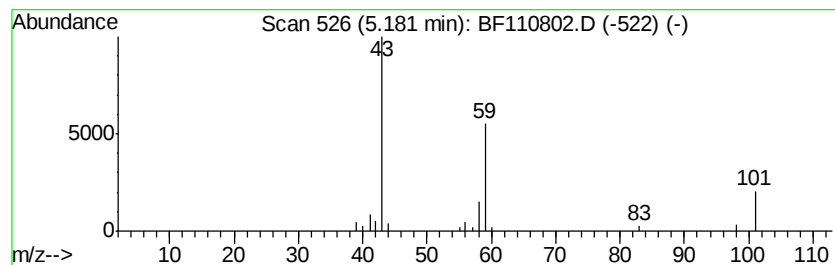
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.38 ng	76616	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Propylamine	59	C3H9N	000107-10-8	9
5		3-Hexanol	102	C6H14O	000623-37-0	9



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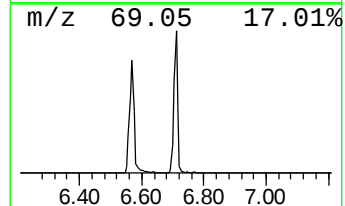
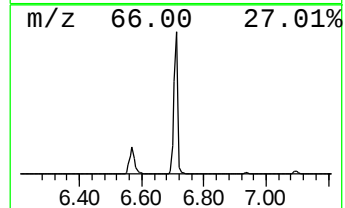
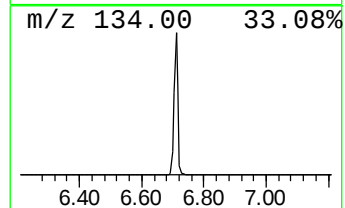
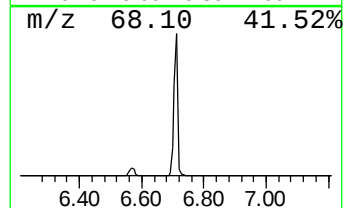
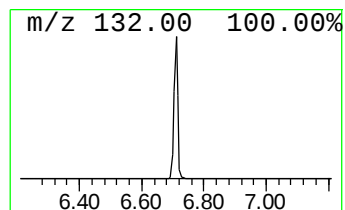
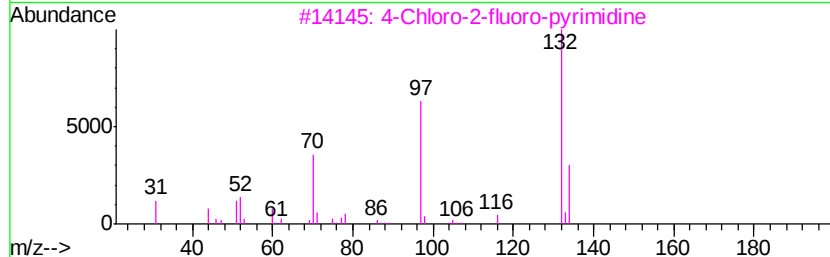
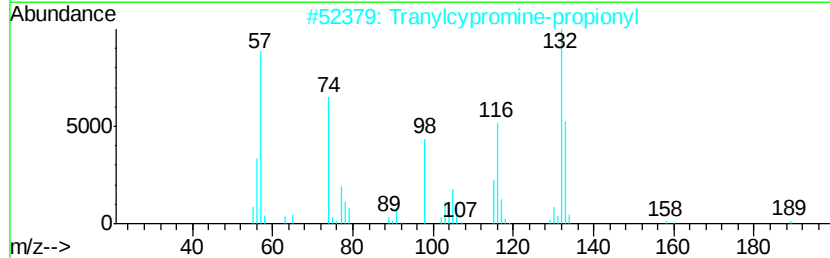
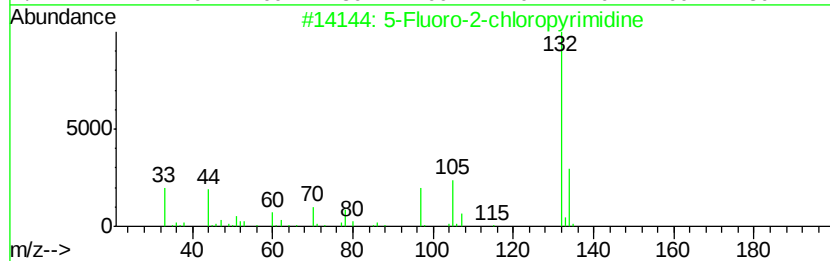
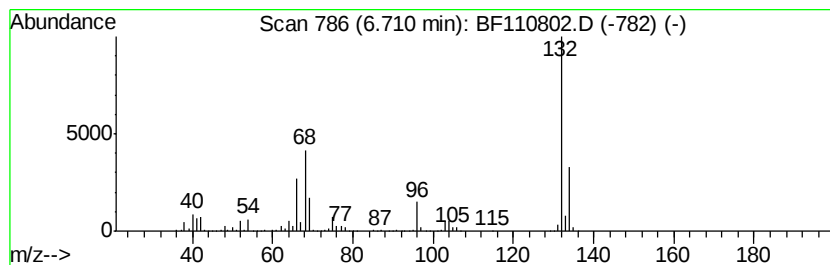
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.71 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	105.07 ng	1839310	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		Xenon	132	Xe	007440-63-3	9



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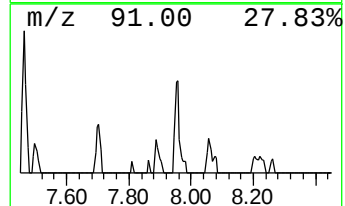
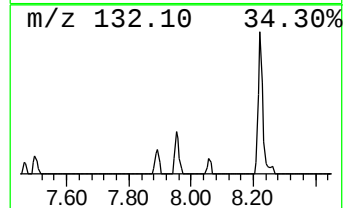
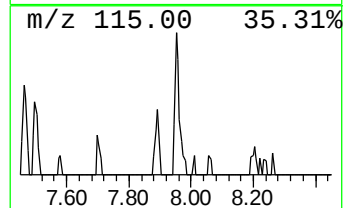
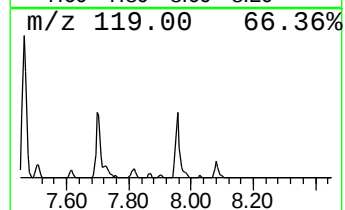
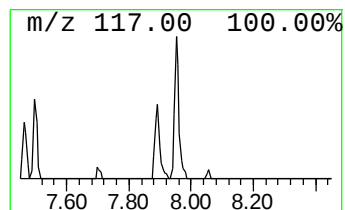
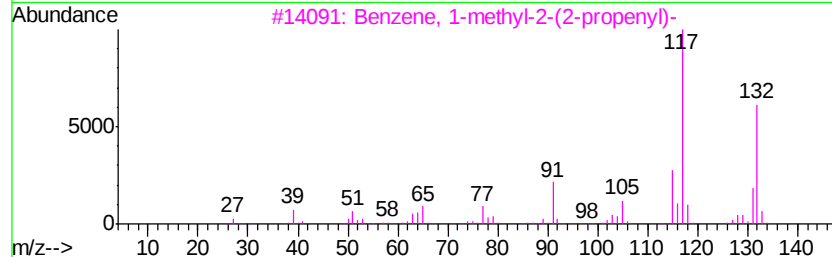
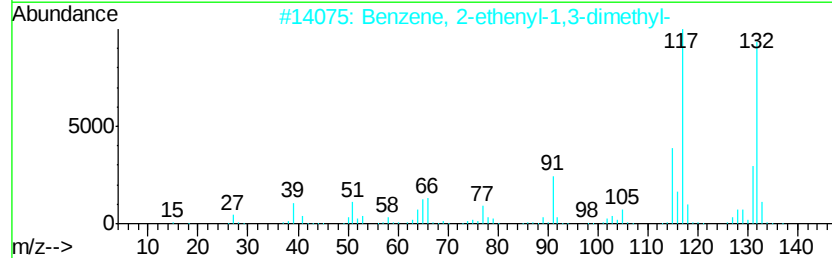
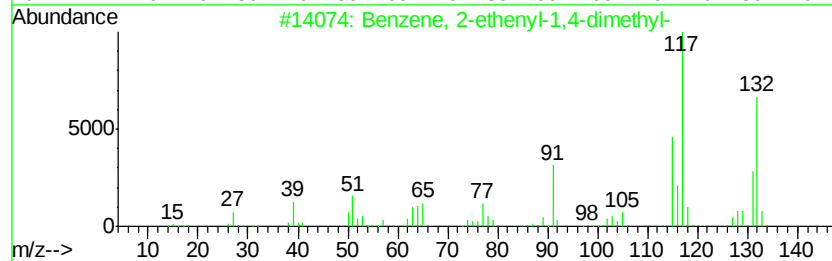
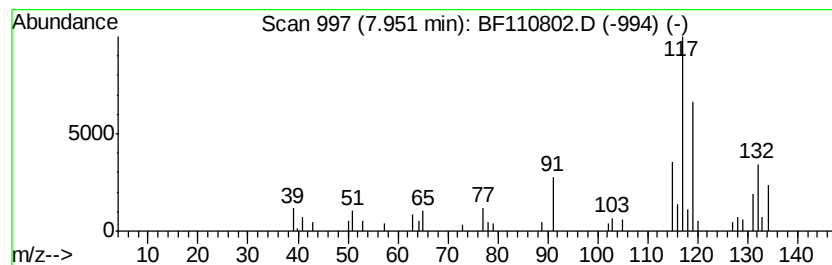
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	2.18 ng	44942	Naphthalene-d8	8.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64



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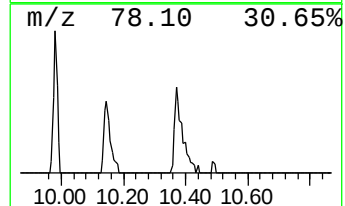
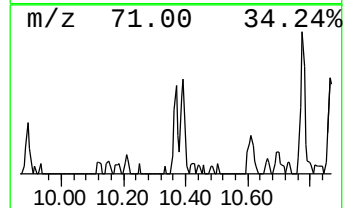
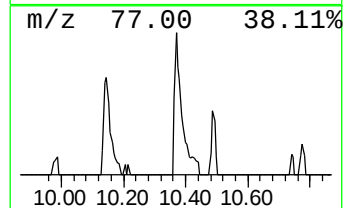
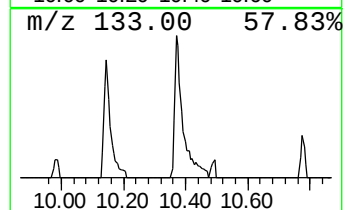
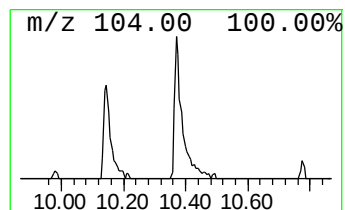
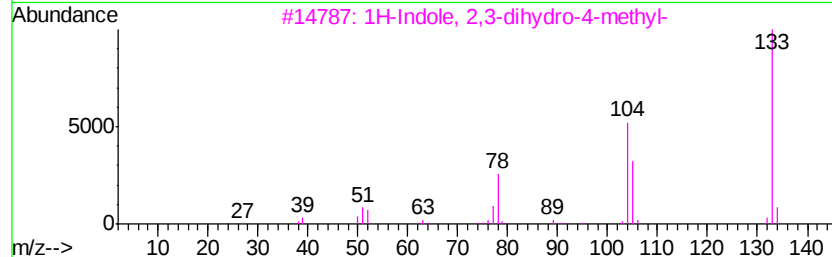
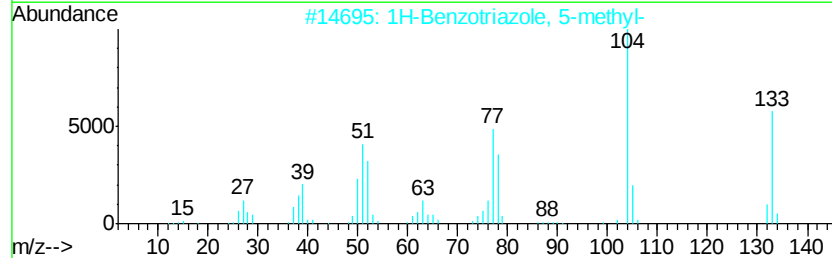
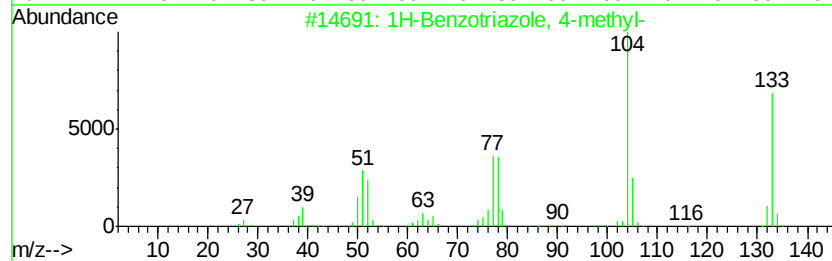
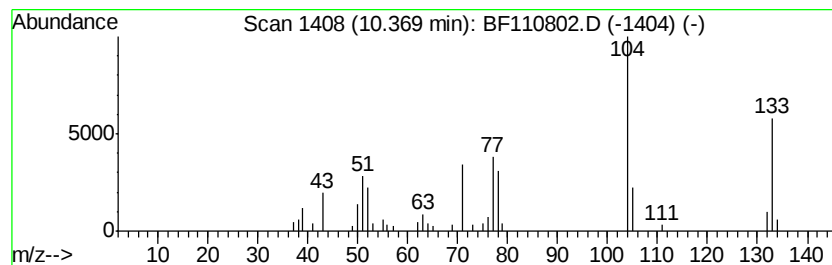
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1H-Benzotriazole, 4-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	2.46 ng	54447	Acenaphthene-d10	9.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	97
2			1H-Benzotriazole, 5-methyl-	133	C7H7N3	000136-85-6	96
3			1H-Indole, 2,3-dihydro-4-methyl-	133	C9H11N	062108-16-1	53
4			1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	38
5			2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
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Operator : JU/SJ
Sample : J5950-06
Misc :
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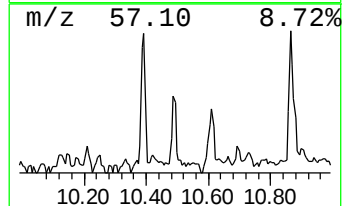
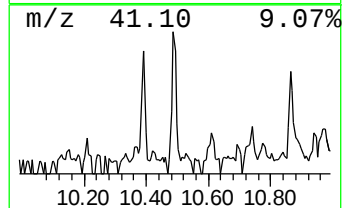
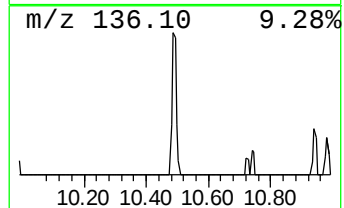
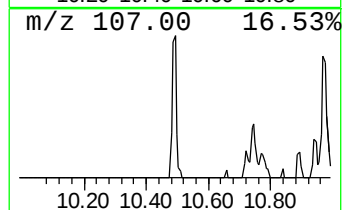
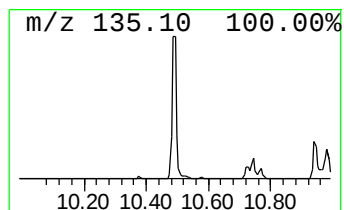
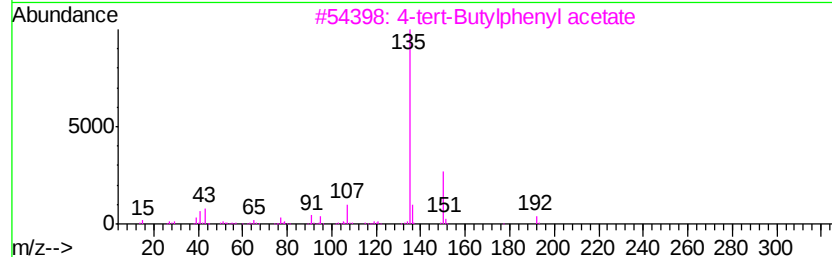
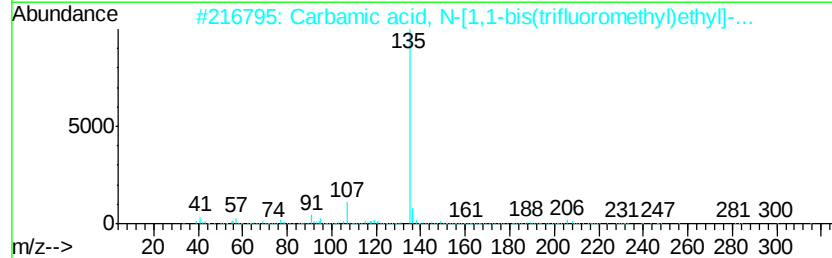
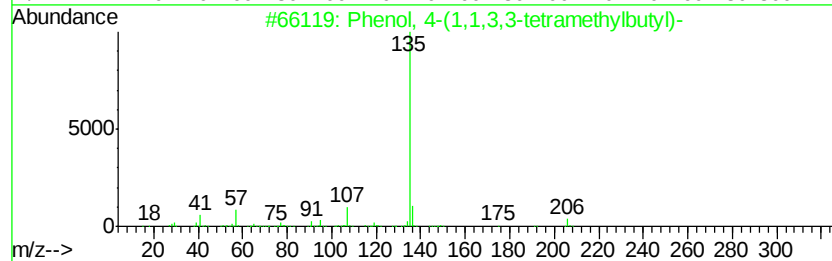
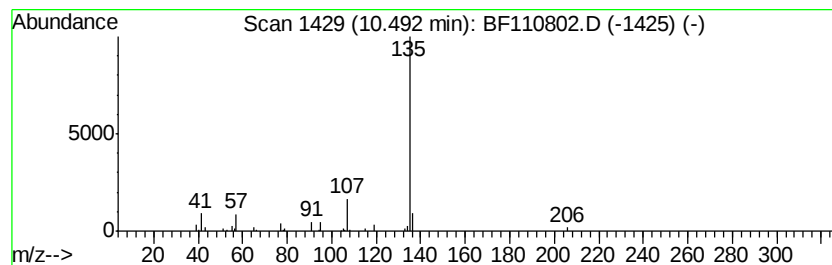
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	3.16 ng	69712	Acenaphthene-d10	9.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206 C14H22O	000140-66-9	83
2	Carbamic acid, N-[1,1-bis(triflu...	413 C19H25F6N02	296242-69-8	72
3	4-tert-Butylphenyl acetate	192 C12H16O2	003056-64-2	72
4	Phenol, 4-(1,1-dimethylpropyl)-	164 C11H16O	000080-46-6	64
5	4-(1,1-Dimethylpropyl)phenyl ace...	206 C13H18O2	1000373-45-3	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
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Sample : J5950-06
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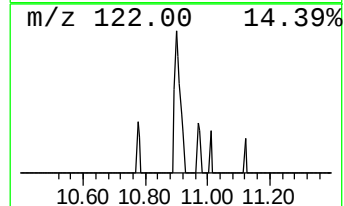
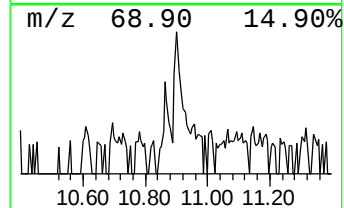
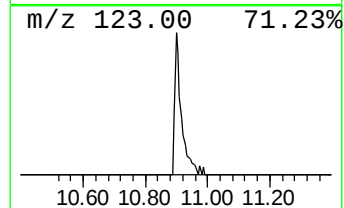
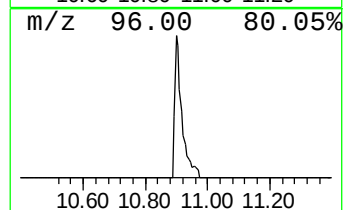
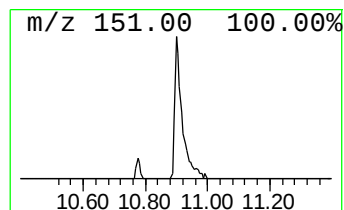
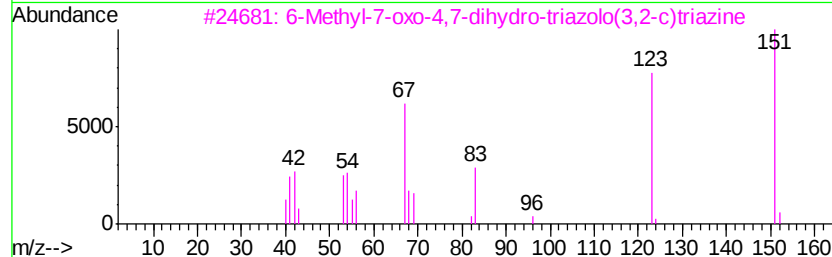
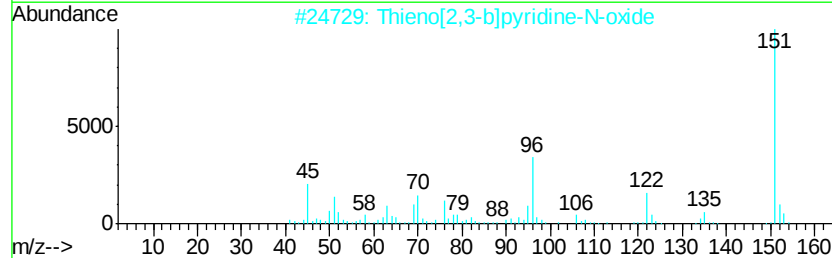
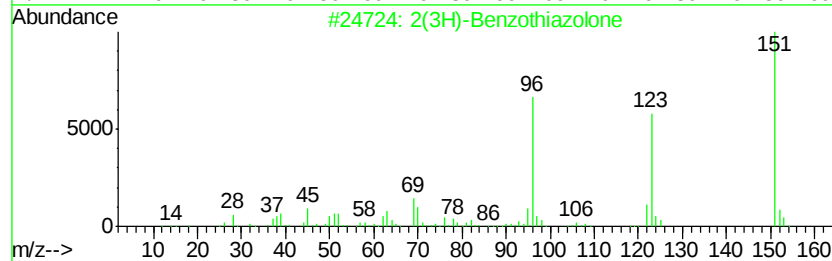
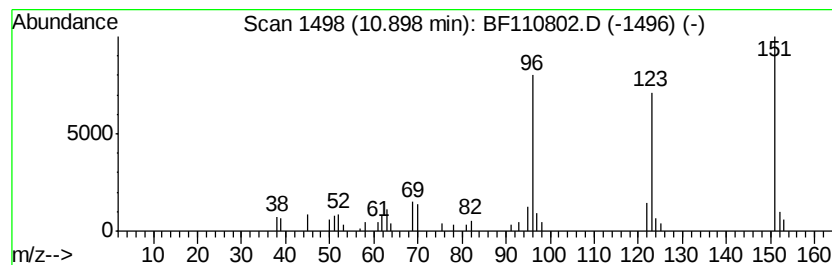
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2(3H)-Benzothiazolone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.90	2.80 ng	49891	Phenanthrene-d10	11.47	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	97
2	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	43
3	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	38
4	4-Methoxyformanilide	151	C8H9NO2	005470-34-8	35
5	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
Data File : BF110802.D
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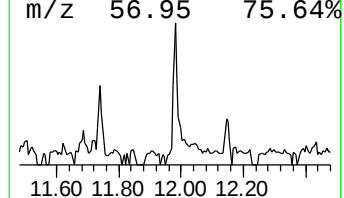
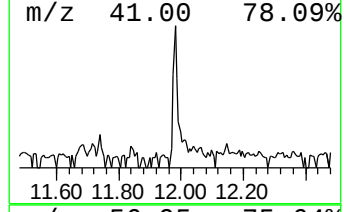
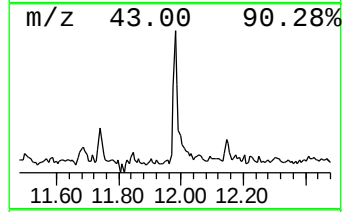
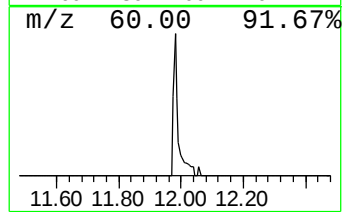
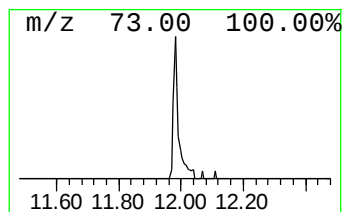
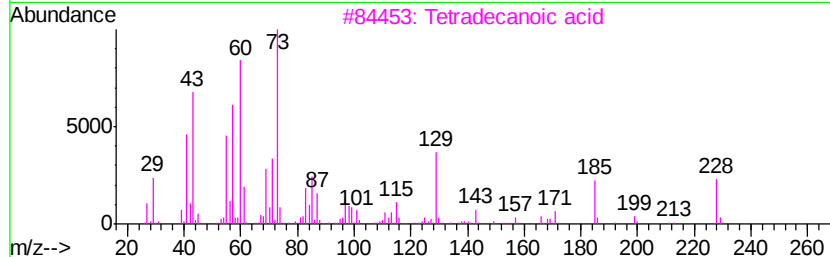
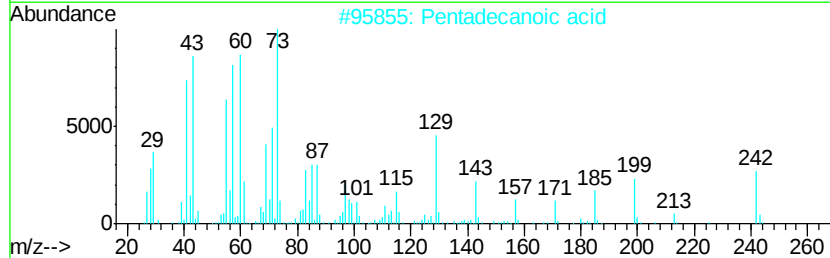
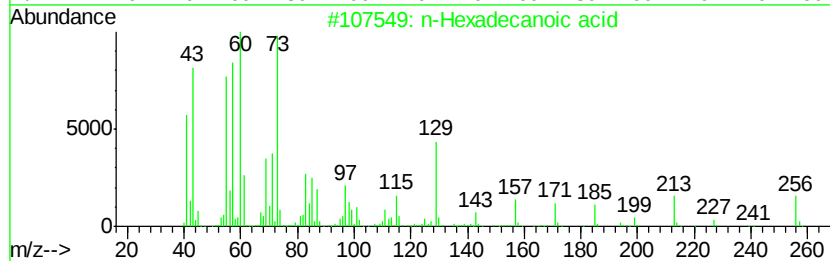
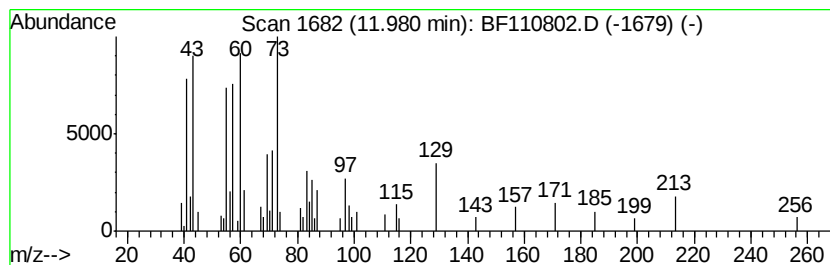
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	3.18 ng	56710	Phenanthrene-d10	11.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4			Tridecanoic acid	214	C13H26O2	000638-53-9	64
5			n-Decanoic acid	172	C10H20O2	000334-48-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
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Acq On : 15 Nov 2018 21:35
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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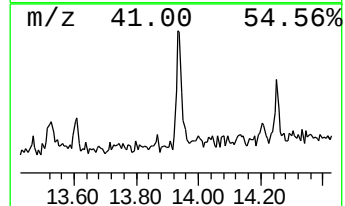
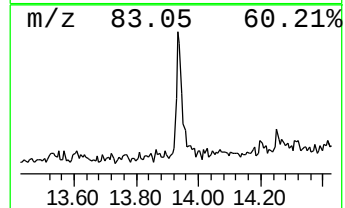
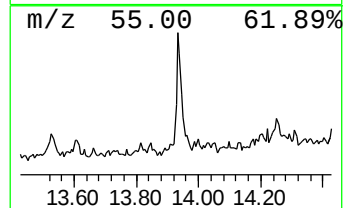
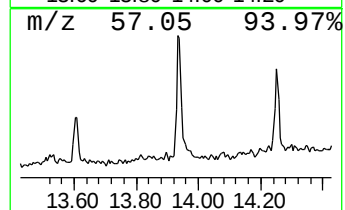
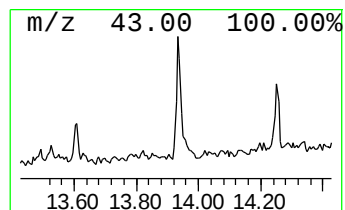
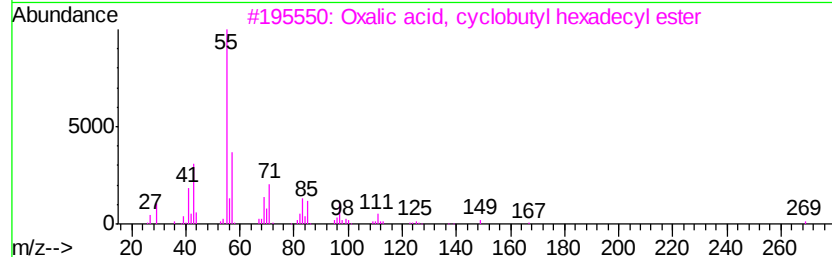
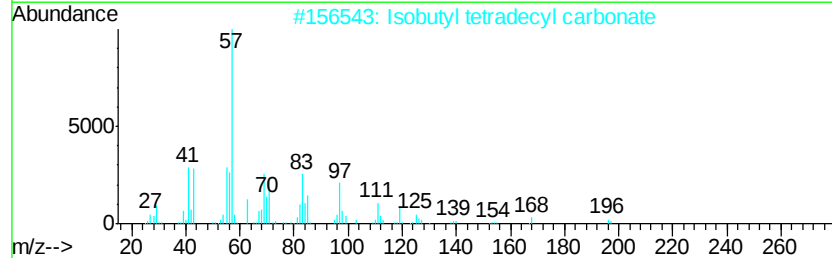
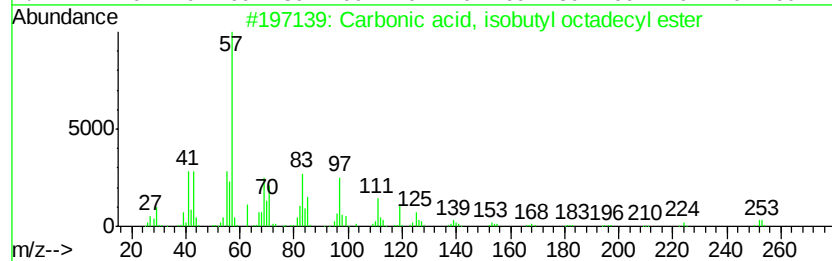
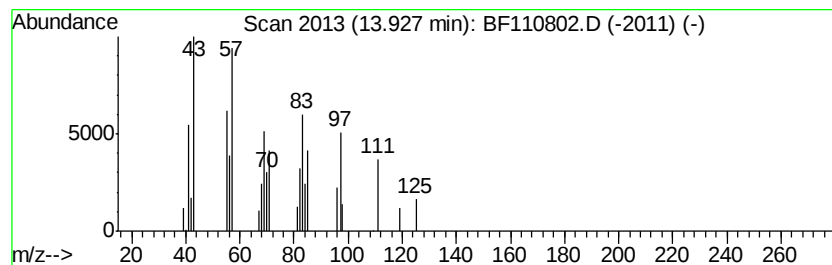
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Carbonic acid, isobutyl oct... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	3.75 ng	68460	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonic acid, isobutyl octadecy...	370	C23H46O3	1000314-61-5	91
2		Isobutyl tetradecyl carbonate	314	C19H38O3	959275-58-2	90
3		Oxalic acid, cyclobutyl hexadecy...	368	C22H40O4	1000309-70-6	90
4		17-Pentatriacontene	491	C35H70	006971-40-0	76
5		1-Docosene	308	C22H44	001599-67-3	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
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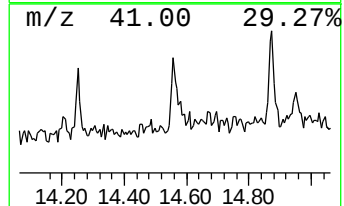
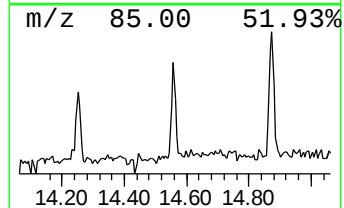
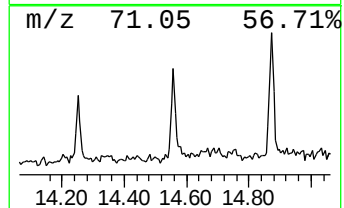
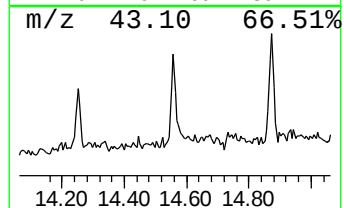
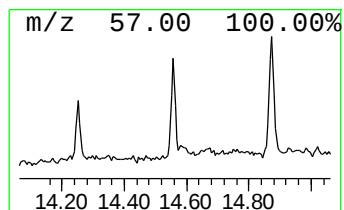
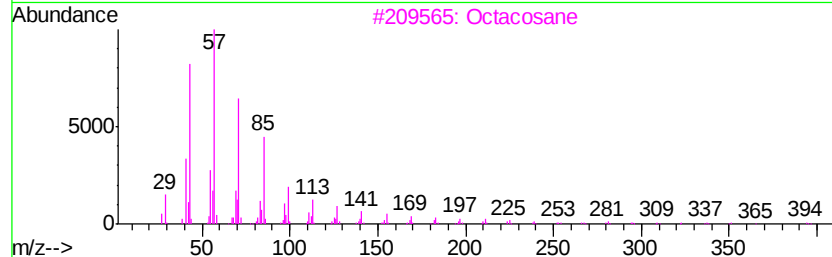
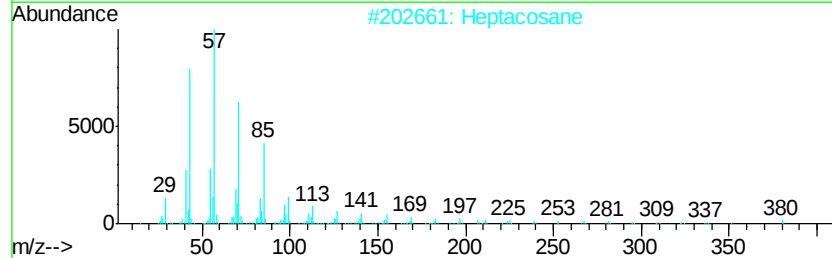
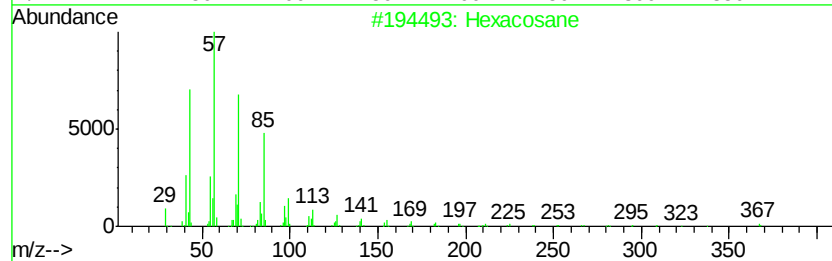
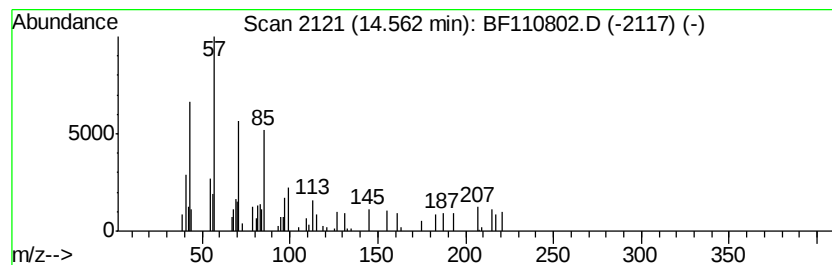
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	3.00 ng	54781	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	90
2		Heptacosane	380	C27H56	000593-49-7	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
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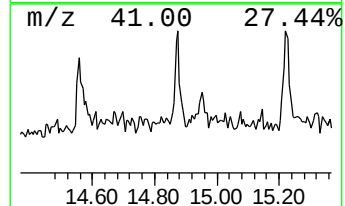
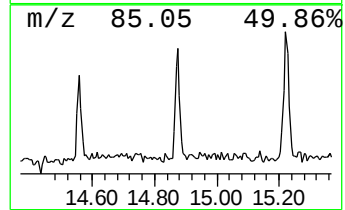
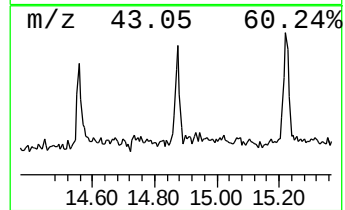
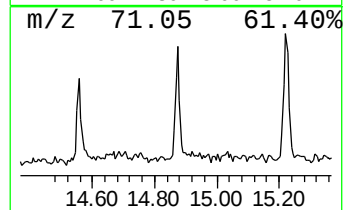
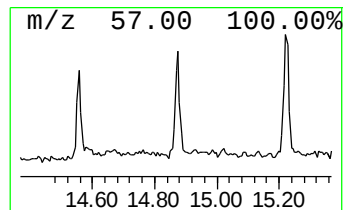
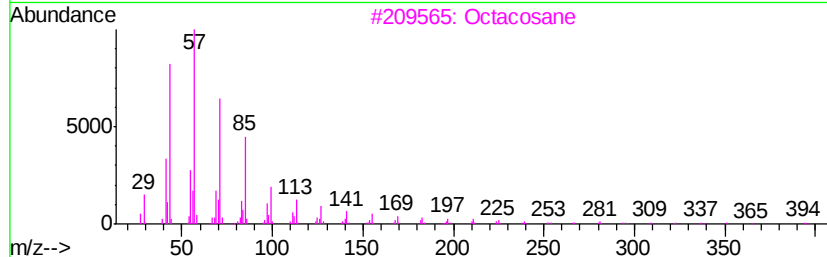
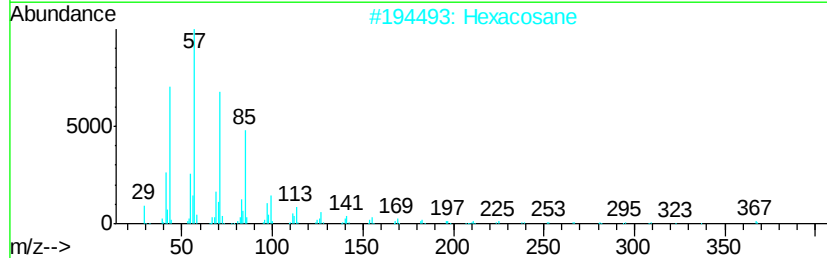
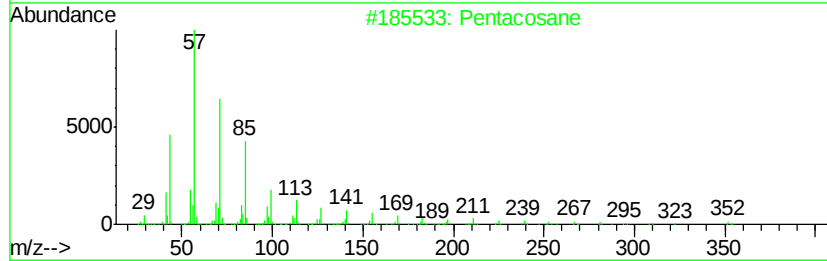
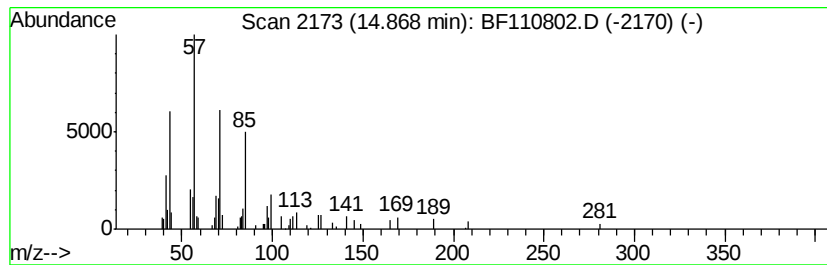
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	2.61 ng	47596	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Octadecane	254	C18H38	000593-45-3	87
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
Data File : BF110802.D
Acq On : 15 Nov 2018 21:35
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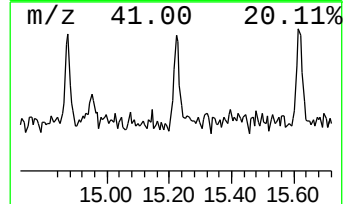
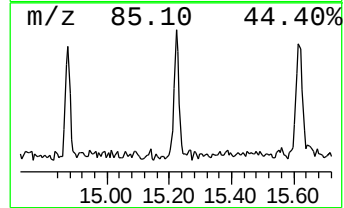
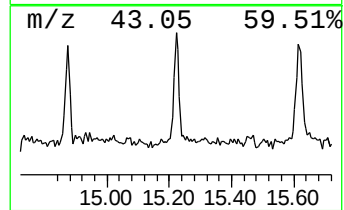
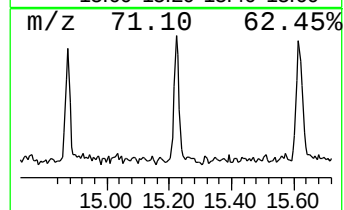
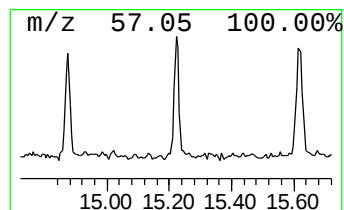
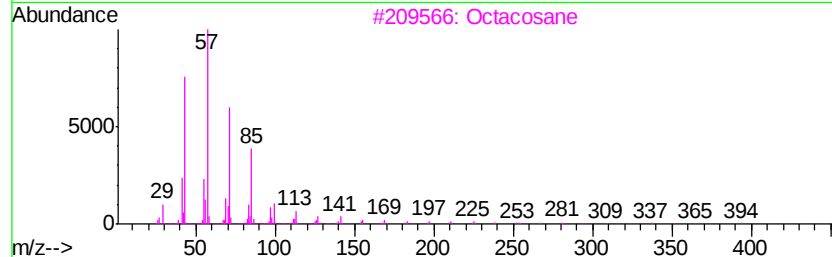
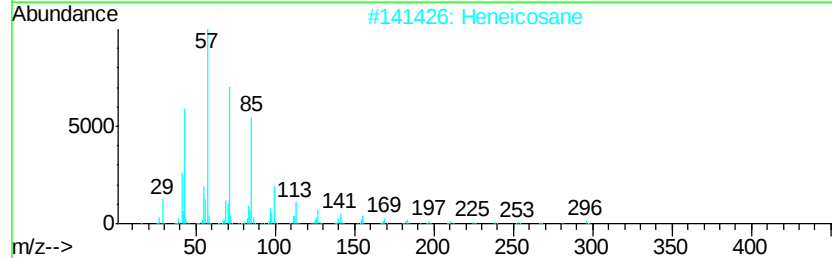
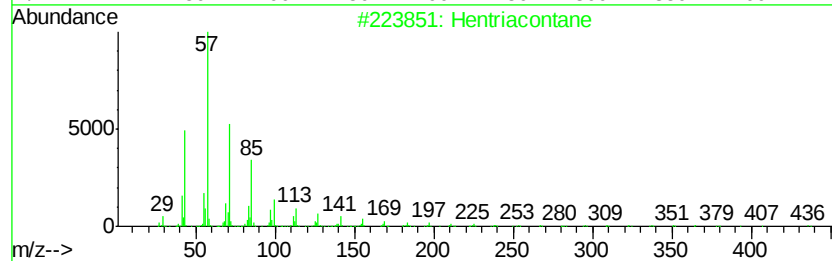
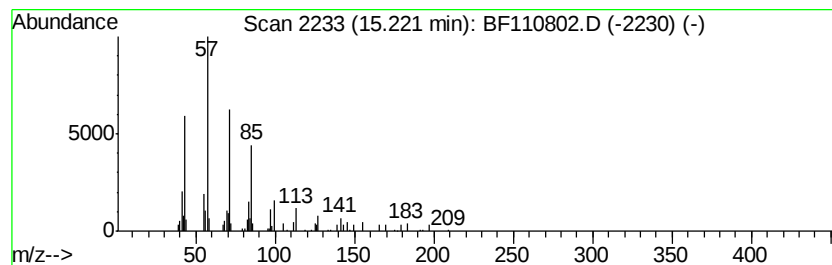
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	3.98 ng	54193	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			Heneicosane	296	C21H44	000629-94-7	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Triacontane	422	C30H62	000638-68-6	87
5			Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
Data File : BF110802.D
Acq On : 15 Nov 2018 21:35
Operator : JU/SJ
Sample : J5950-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSI-20-TW-1

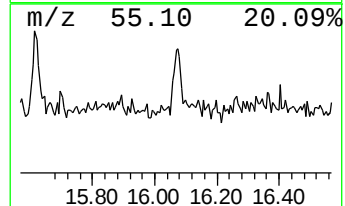
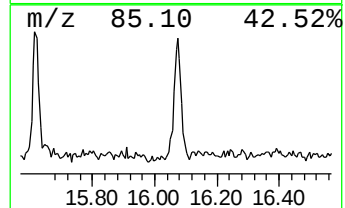
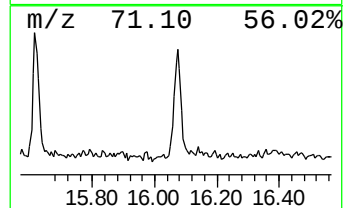
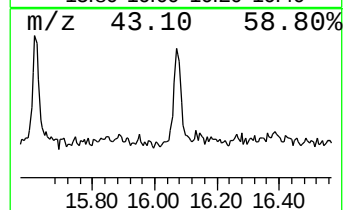
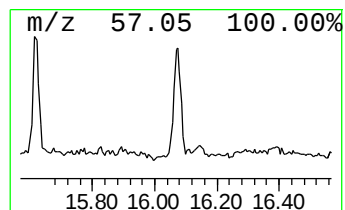
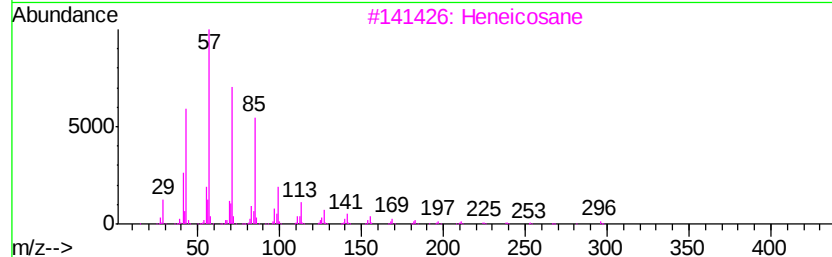
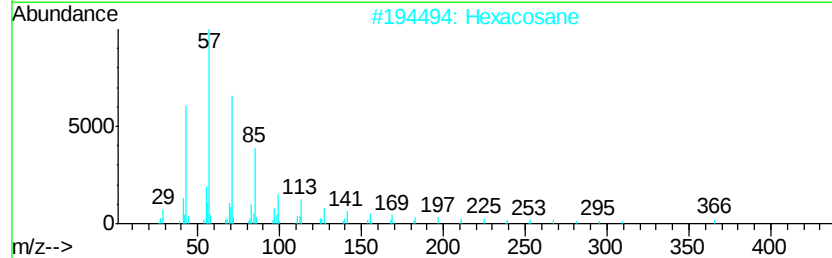
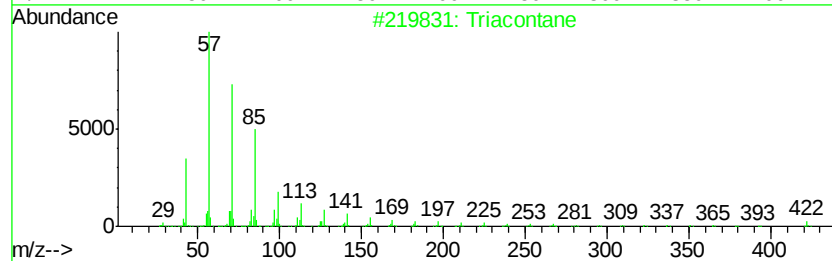
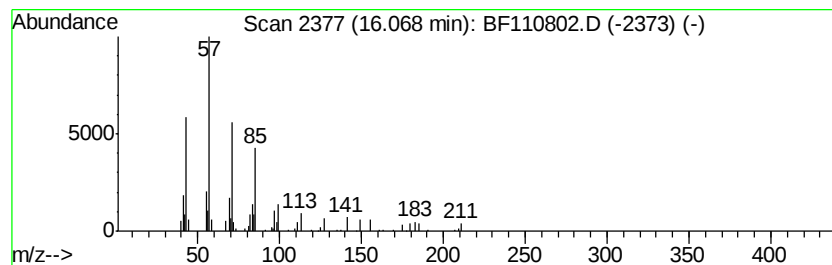
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Triacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	4.74 ng	64548	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			Hexacosane	366	C26H54	000630-01-3	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Pentacosane	352	C25H52	000629-99-2	90
5			Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
Data File : BF110802.D
Acq On : 15 Nov 2018 21:35
Operator : JU/SJ
Sample : J5950-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSI-20-TW-1

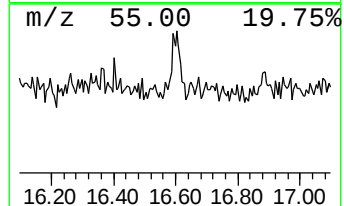
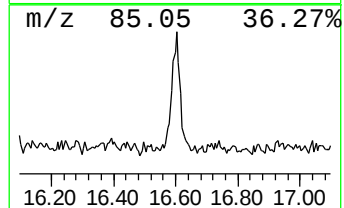
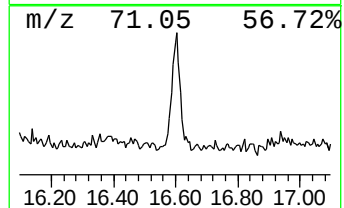
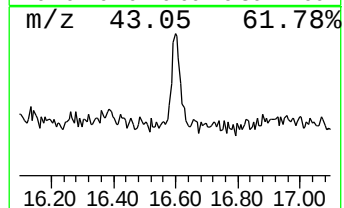
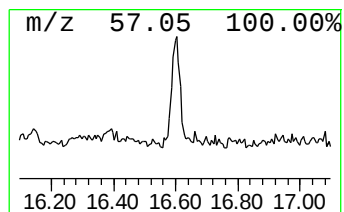
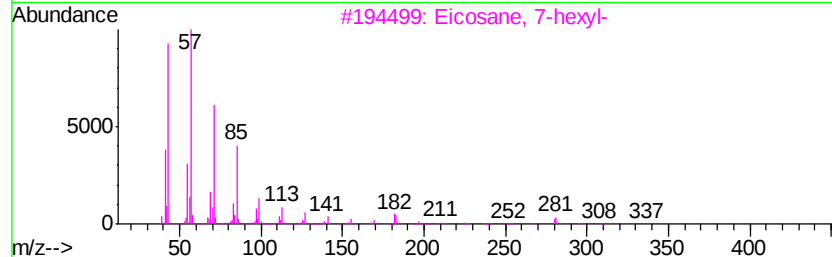
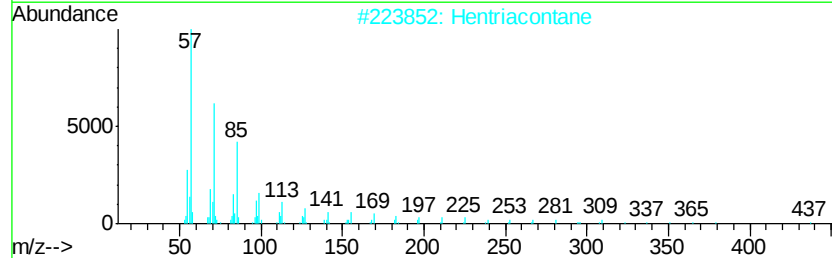
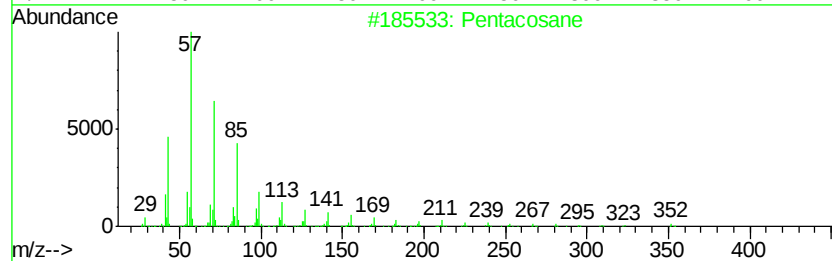
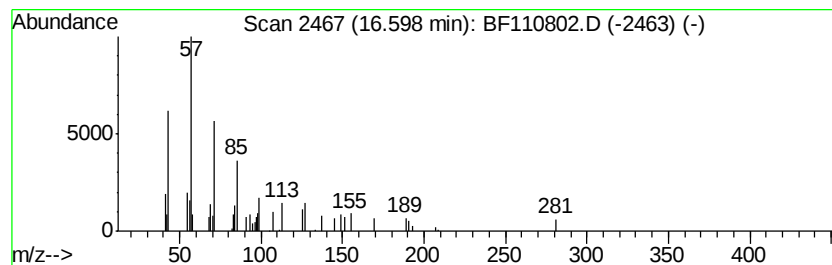
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexatriacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.60	3.32 ng	45168	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	86
2		Hentriacontane	437	C31H64	000630-04-6	86
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	83
4		Hexatriacontane	507	C36H74	000630-06-8	78
5		Heptacosane	380	C27H56	000593-49-7	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
Data File : BF110802.D
Acq On : 15 Nov 2018 21:35
Operator : JU/SJ
Sample : J5950-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSI-20-TW-1

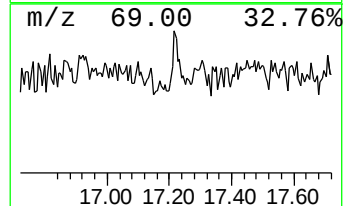
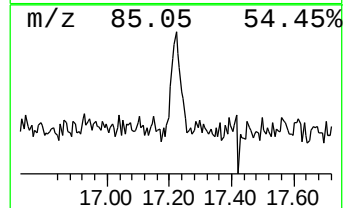
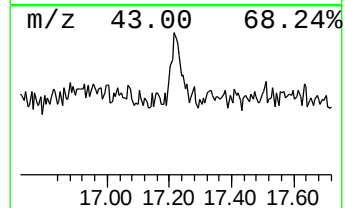
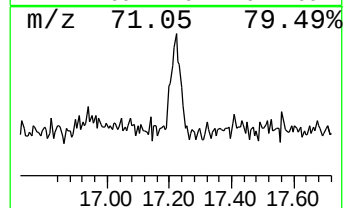
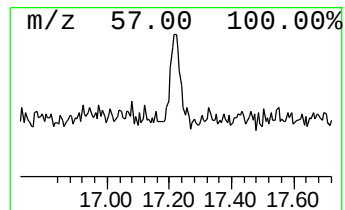
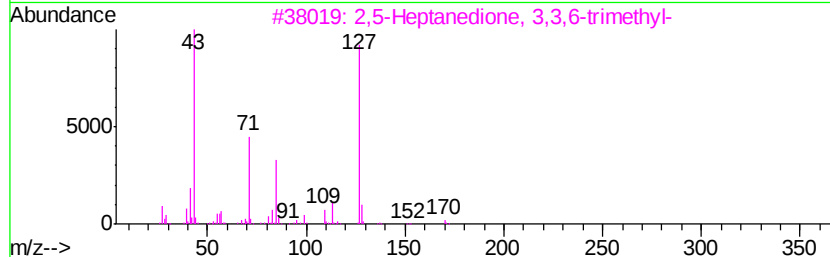
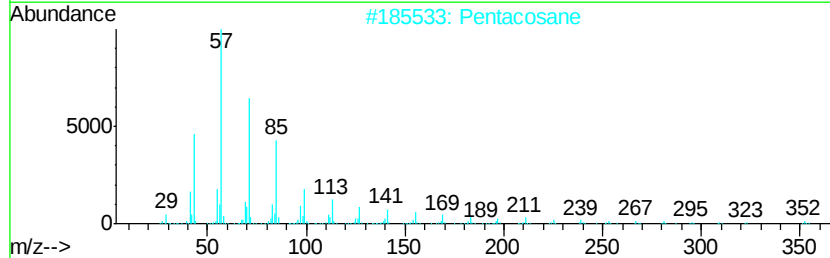
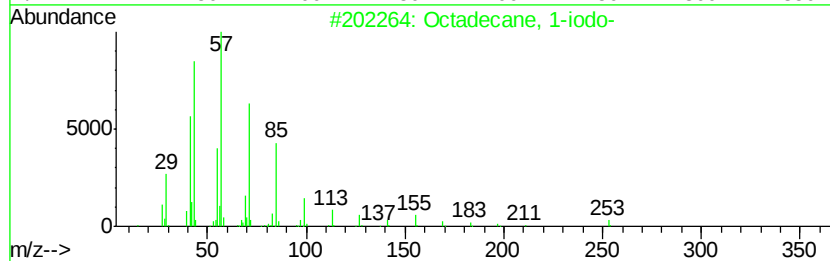
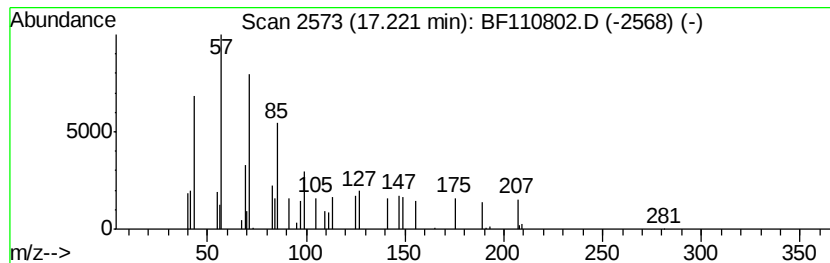
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Octadecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.22	2.20 ng	29944	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	53
2		Pentacosane	352	C25H52	000629-99-2	53
3		2,5-Heptanedione, 3,3,6-trimethyl-	170	C10H18O2	051513-40-7	50
4		Triacontane	422	C30H62	000638-68-6	47
5		Heneicosane	296	C21H44	000629-94-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
 Data File : BF110802.D
 Acq On : 15 Nov 2018 21:35
 Operator : JU/SJ
 Sample : J5950-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.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...	2.24	8.6	ng	150990	1	6.94	350103	20.0
2-Pentanone, 4-hy...	5.18	4.4	ng	76616	1	6.94	350103	20.0
unknown6.71	6.71	105.1	ng	1839310	1	6.94	350103	20.0
Benzene, 2-etheny...	7.95	2.2	ng	44942	2	8.22	412817	20.0
1H-Benzotriazole,...	10.37	2.5	ng	54447	3	9.98	441903	20.0
Phenol, 4-(1,1,3,...	10.49	3.2	ng	69712	3	9.98	441903	20.0
2(3H)-Benzothiazo...	10.90	2.8	ng	49891	4	11.47	356230	20.0
n-Hexadecanoic acid	11.98	3.2	ng	56710	4	11.47	356230	20.0
Carbonic acid, is...	13.93	3.8	ng	68460	5	14.13	365351	20.0
Hexacosane	14.56	3.0	ng	54781	5	14.13	365351	20.0
Pentacosane	14.87	2.6	ng	47596	5	14.13	365351	20.0
Hentriacontane	15.22	4.0	ng	54193	6	15.66	272228	20.0
Triacontane	16.07	4.7	ng	64548	6	15.66	272228	20.0
Hexatriacontane	16.60	3.3	ng	45168	6	15.66	272228	20.0
Octadecane, 1-iodo-	17.22	2.2	ng	29944	6	15.66	272228	20.0