

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
 Data File : BF103863.D
 Acq On : 22 Mar 2018 20:18
 Operator : JU/SJ
 Sample : J1997-25
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SLAB-CONCRETE-B

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-BF031518.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	2.304	32	37	43	rVB	113741	146714	2.22%	0.441%
2	3.787	282	289	295	rBV	86430	121731	1.84%	0.366%
3	4.628	422	432	435	rBV	3323135	5140531	77.82%	15.464%
4	5.240	525	536	539	rBV	3048741	6606058	100.00%	19.872%
5	5.592	592	596	610	rVB	560278	484772	7.34%	1.458%
6	6.592	761	766	769	rBV	1637355	1579804	23.91%	4.752%
7	6.739	788	791	794	rBV	934814	741675	11.23%	2.231%
8	6.869	809	813	820	rBV2	69673	86822	1.31%	0.261%
9	6.975	827	831	834	rVB	1017894	857684	12.98%	2.580%
10	7.128	853	857	861	rVB	2513223	2364234	35.79%	7.112%
11	7.292	881	885	889	rBV3	111803	116221	1.76%	0.350%
12	7.534	922	926	929	rBV	1977451	1801022	27.26%	5.418%
13	7.616	934	940	943	rBV5	88108	131455	1.99%	0.395%
14	7.657	943	947	952	rBV6	34278	67940	1.03%	0.204%
15	7.757	962	964	966	rVV2	82433	68543	1.04%	0.206%
16	7.845	975	979	984	rBV3	50555	90472	1.37%	0.272%
17	7.945	993	996	1000	rVB	438169	347623	5.26%	1.046%
18	8.257	1045	1049	1054	rBV	1182560	1071269	16.22%	3.223%
19	9.192	1205	1208	1212	rBV	685302	548036	8.30%	1.649%
20	9.310	1224	1228	1229	rBV	979554	953074	14.43%	2.867%
21	9.328	1229	1231	1234	rVB	3567740	2878508	43.57%	8.659%
22	9.427	1245	1248	1253	rVB	85179	74224	1.12%	0.223%
23	9.716	1293	1297	1300	rBV	261278	215245	3.26%	0.647%
24	9.927	1329	1333	1337	rBV	124117	98522	1.49%	0.296%
25	10.016	1344	1348	1354	rVB	1104757	1041927	15.77%	3.134%
26	10.404	1411	1414	1416	rBV	119880	104526	1.58%	0.314%
27	10.427	1416	1418	1421	rVB	147964	109794	1.66%	0.330%
28	10.904	1495	1499	1504	rBV	132793	178168	2.70%	0.536%
29	11.351	1572	1575	1577	rBV	95878	70332	1.06%	0.212%
30	11.380	1577	1580	1583	rVB2	99342	83819	1.27%	0.252%
31	11.504	1598	1601	1605	rBV	930614	852283	12.90%	2.564%
32	13.092	1868	1871	1877	rVB	2376053	2327690	35.24%	7.002%
33	13.974	2019	2021	2025	rVB	215903	186219	2.82%	0.560%
34	14.121	2043	2046	2048	rBV	178901	163165	2.47%	0.491%

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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	14.162	2049	2053	2056	rVB	818787	813470	12.31%	2.447%
36	15.704	2310	2315	2321	rVB	582587	719040	10.88%	2.163%

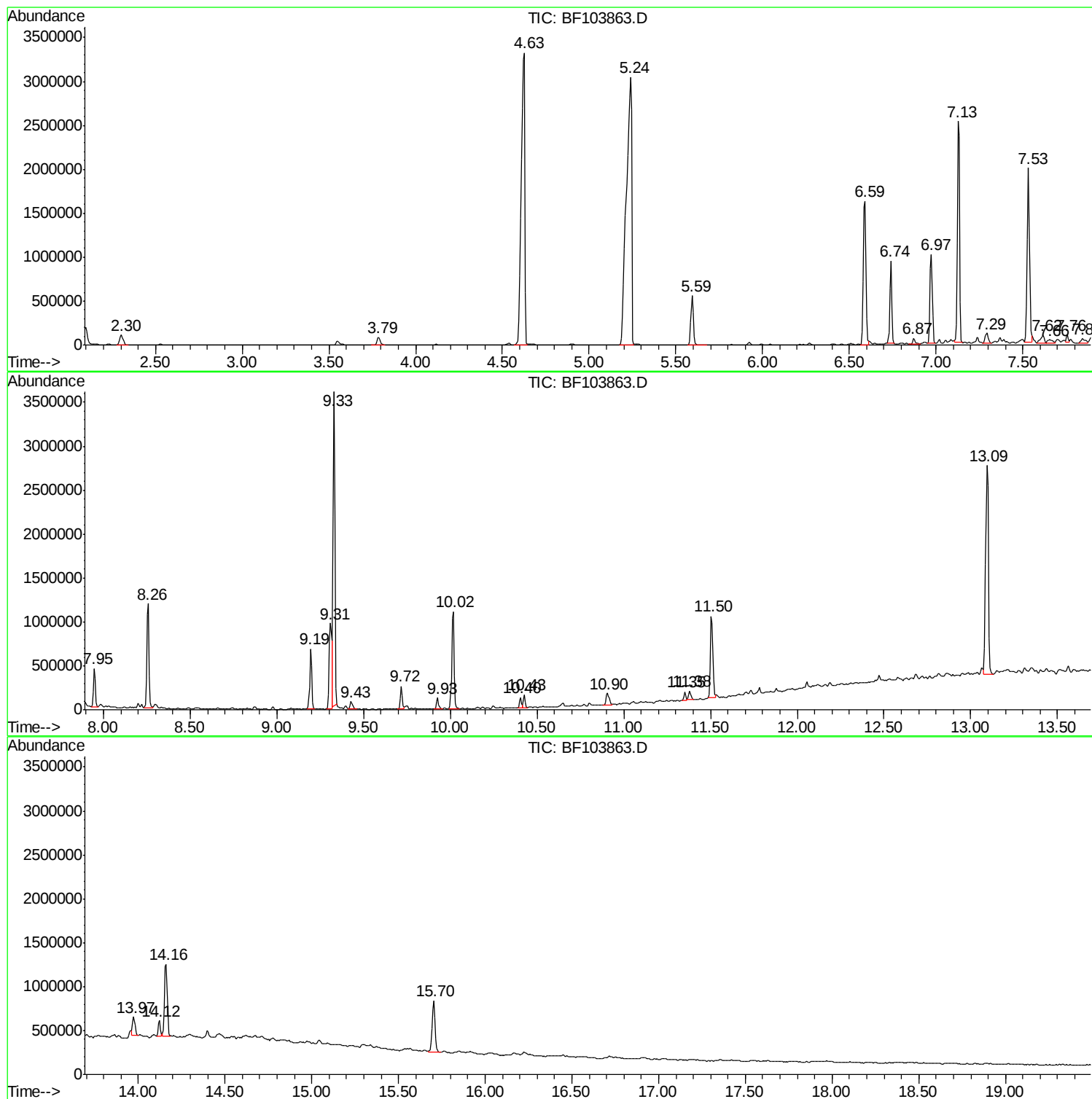
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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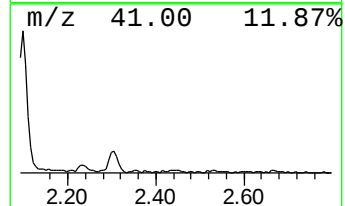
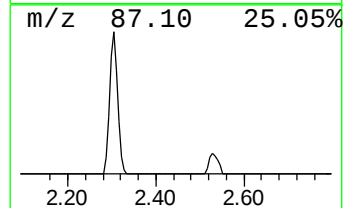
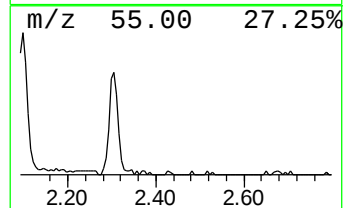
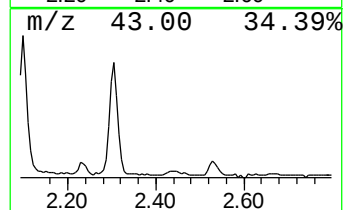
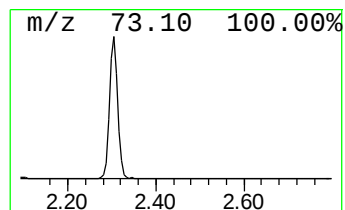
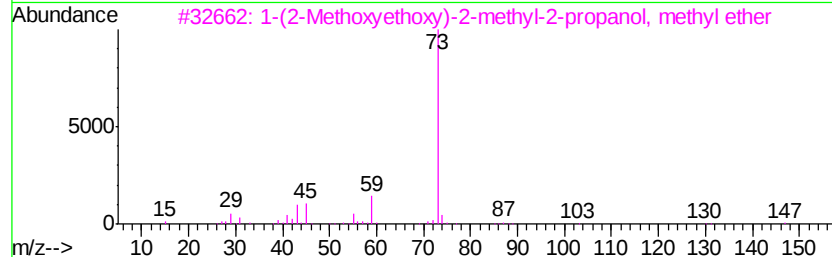
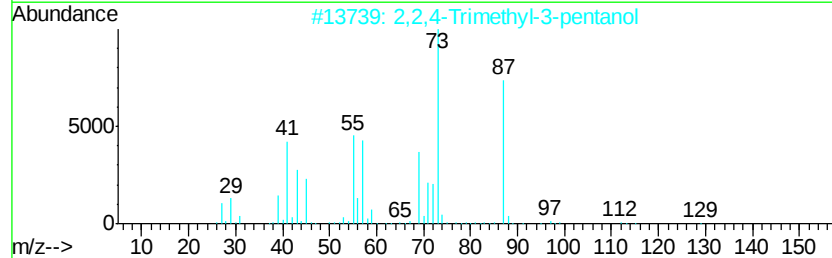
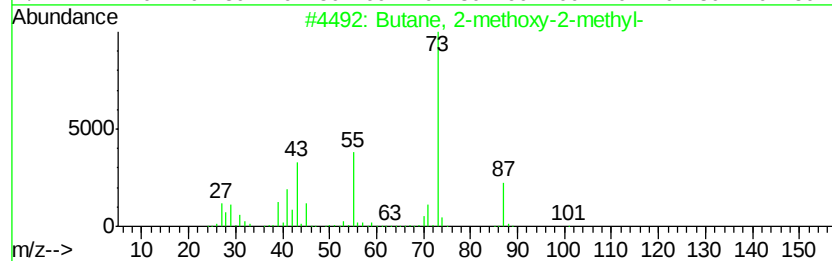
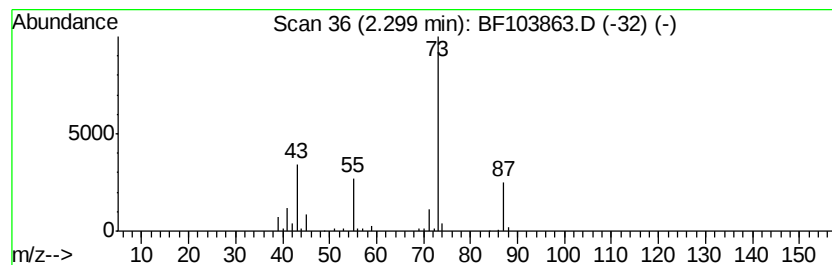
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.30	3.42 ng	146714	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17



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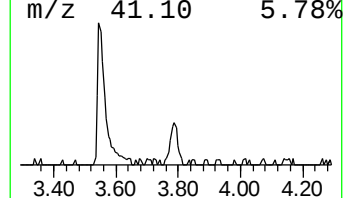
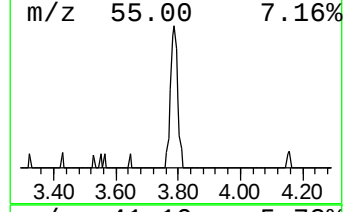
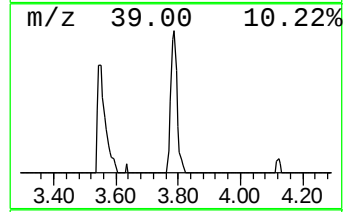
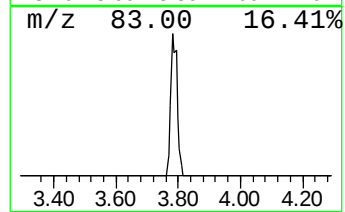
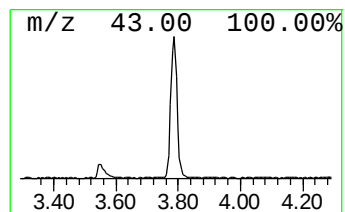
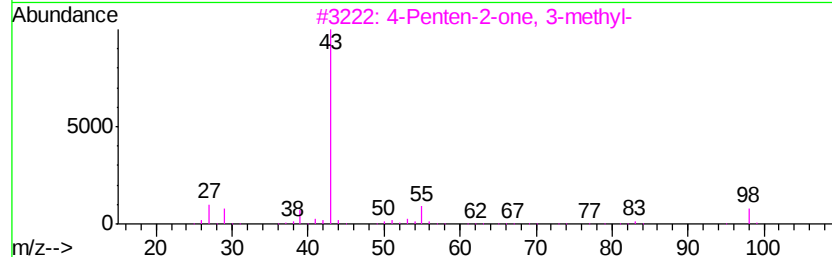
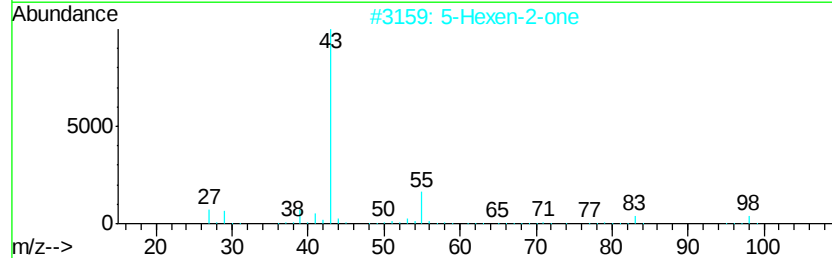
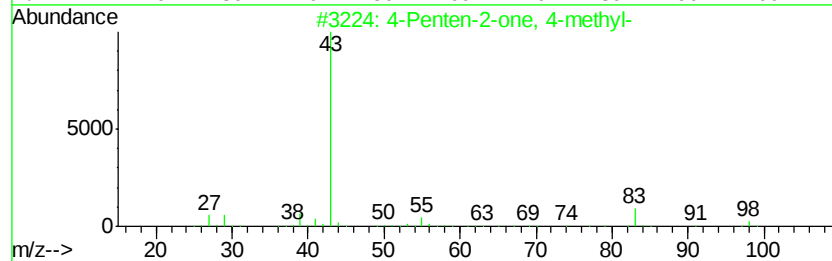
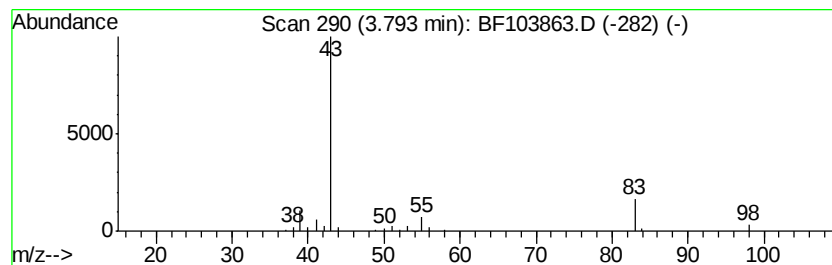
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4-Penten-2-one, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.79	2.84 ng	121731	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	80
2			5-Hexen-2-one	98	C6H10O	000109-49-9	46
3			4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	43
4			2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	12
5			Isoxazole, 5-methyl-	83	C4H5NO	005765-44-6	9



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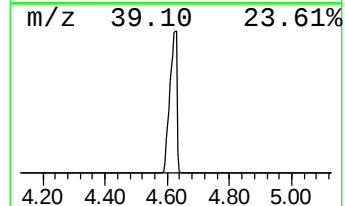
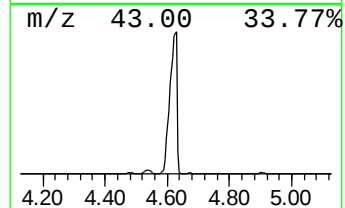
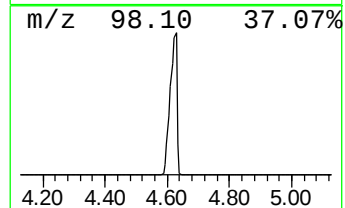
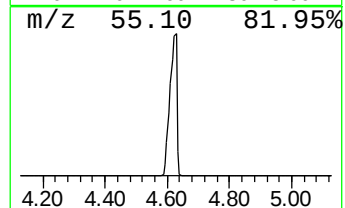
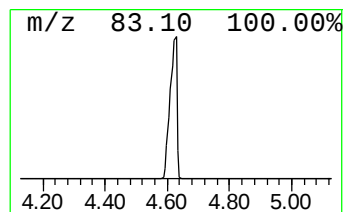
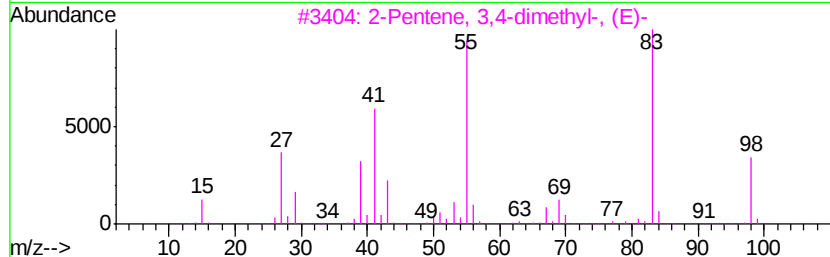
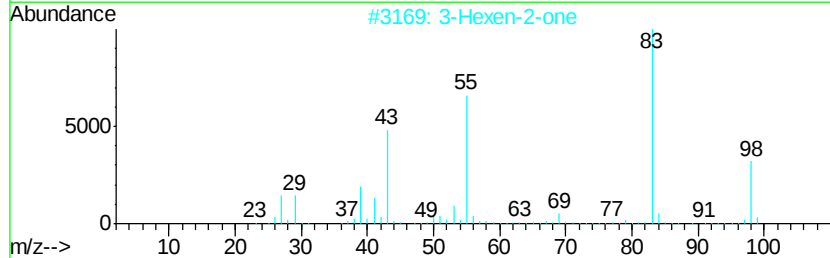
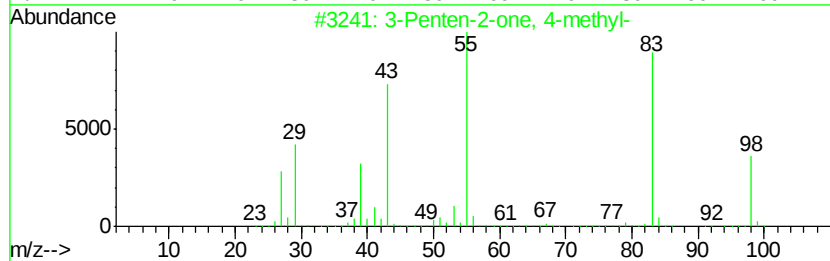
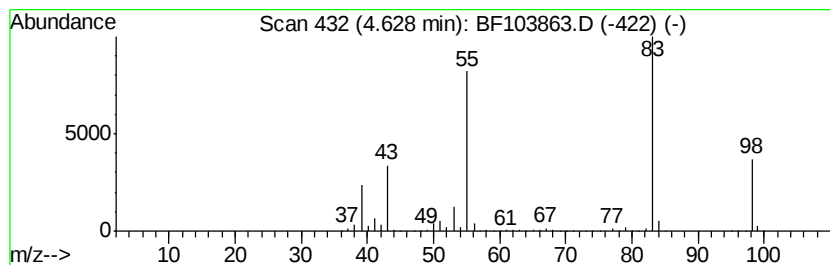
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	119.87 ng	5140530	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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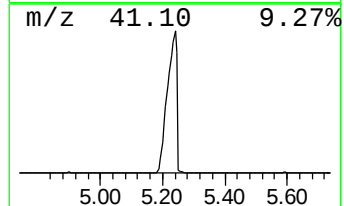
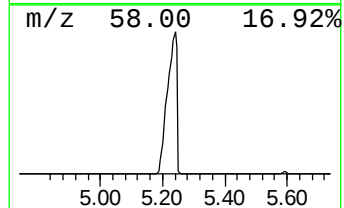
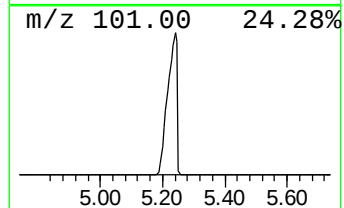
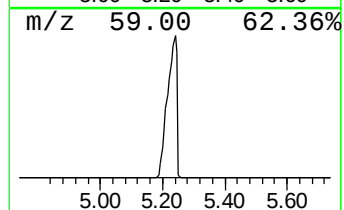
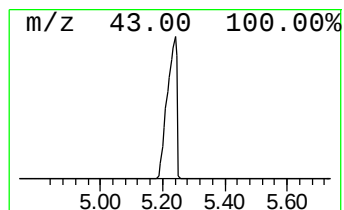
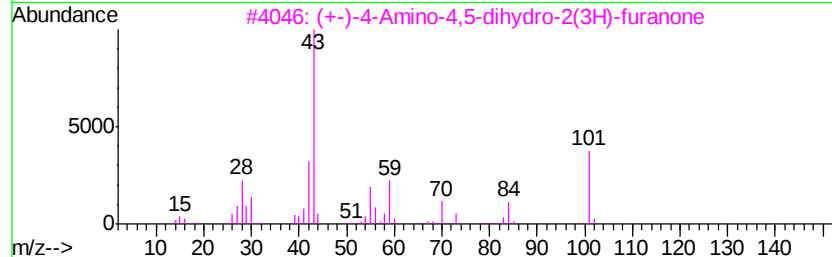
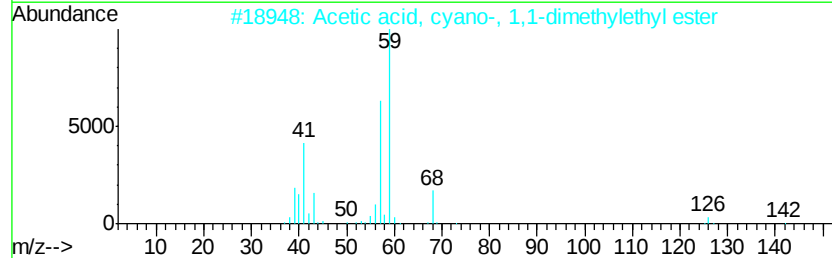
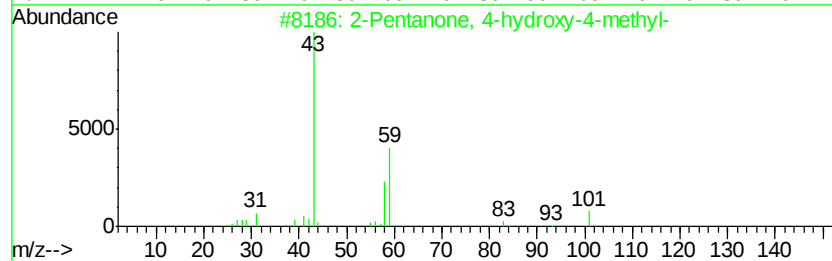
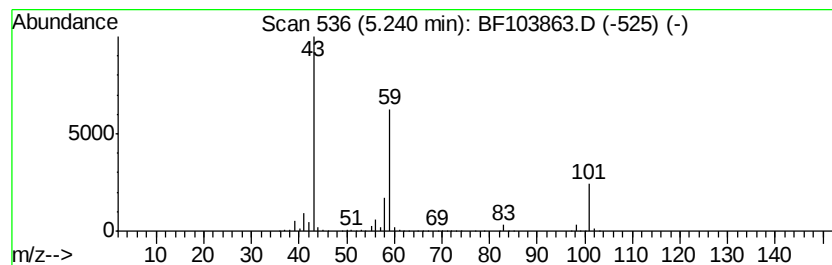
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	154.04 ng	6606060	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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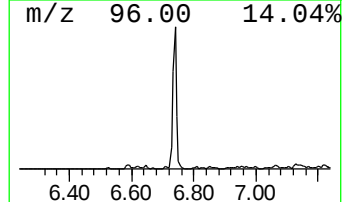
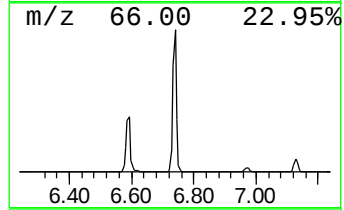
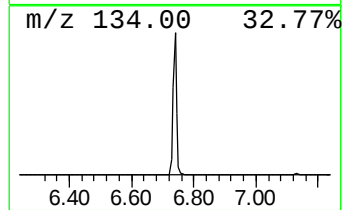
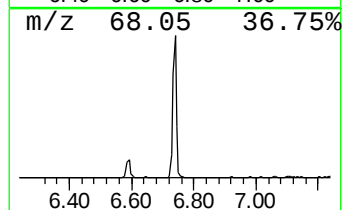
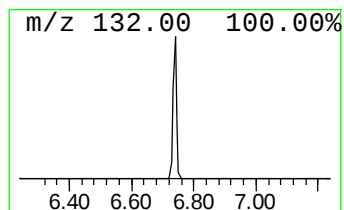
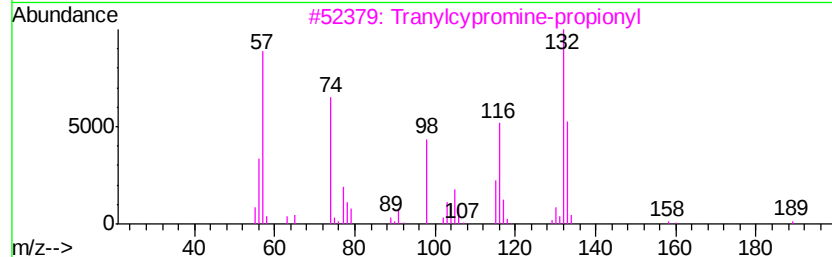
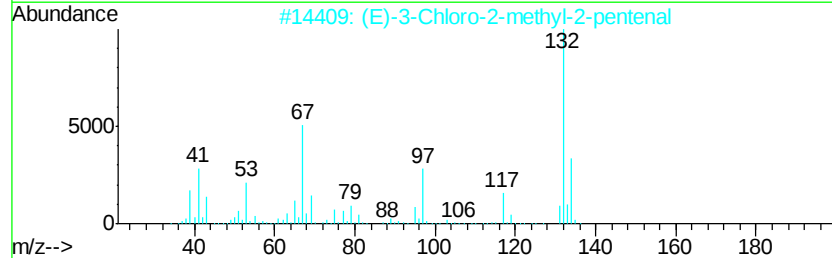
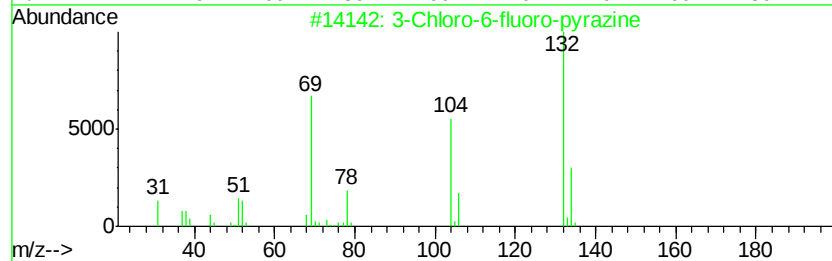
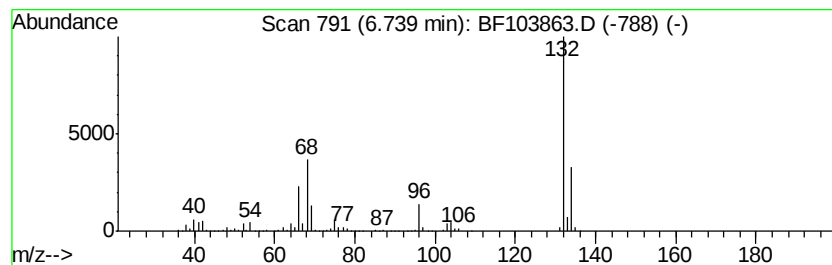
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown6.74 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	17.29 ng	741675	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	12



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Data File : BF103863.D
Acq On : 22 Mar 2018 20:18
Operator : JU/SJ
Sample : J1997-25
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SLAB-CONCRETE-B

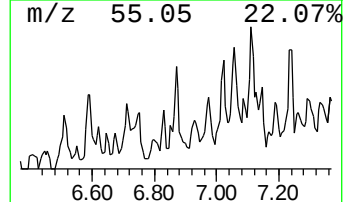
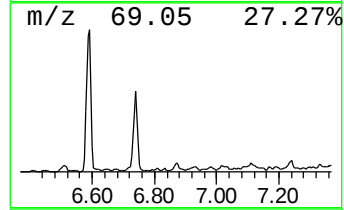
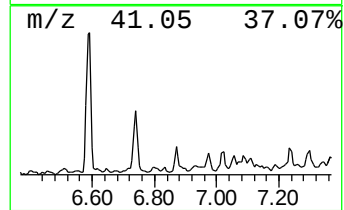
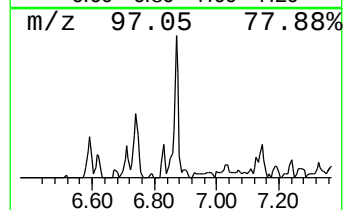
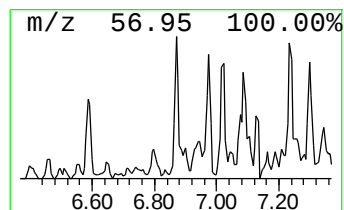
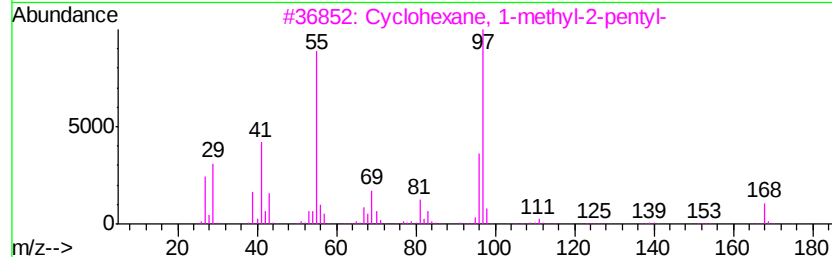
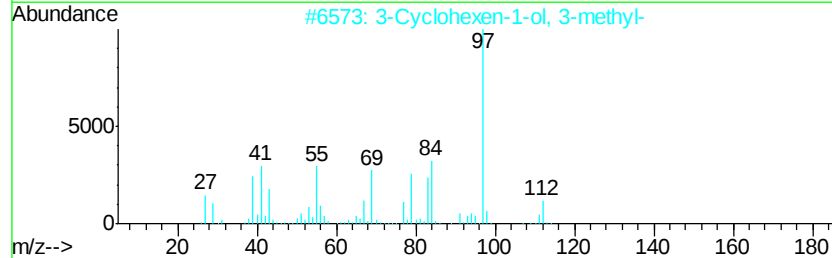
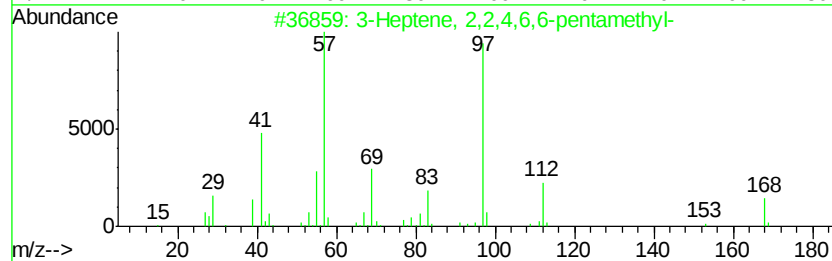
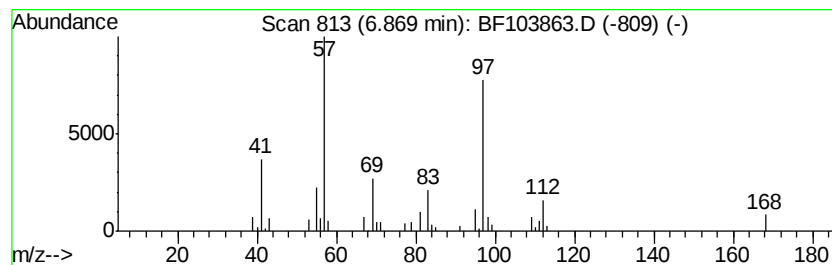
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 3-Heptene, 2,2,4,6,6-pentam... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.87	2.02 ng	86822	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Heptene, 2,2,4,6,6-pentamethyl-	168	C12H24	000123-48-8	86
2			3-Cyclohexen-1-ol, 3-methyl-	112	C7H12O	053783-91-8	58
3			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	50
4			2-Pentene, 2,4,4-trimethyl-	112	C8H16	000107-40-4	49
5			2-Pentene, 3,4,4-trimethyl-	112	C8H16	000598-96-9	47



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF032218\
Data File : BF103863.D
Acq On : 22 Mar 2018 20:18
Operator : JU/SJ
Sample : J1997-25
Misc :
ALS Vial : 25 Sample Multiplier: 1

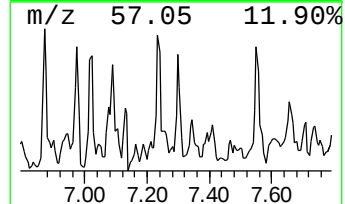
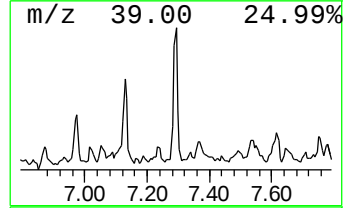
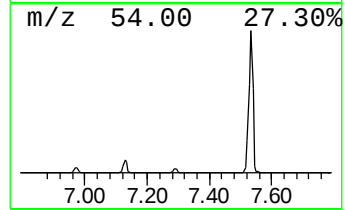
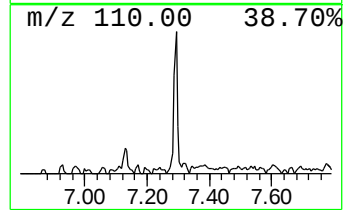
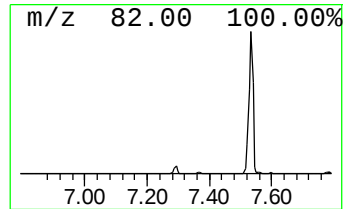
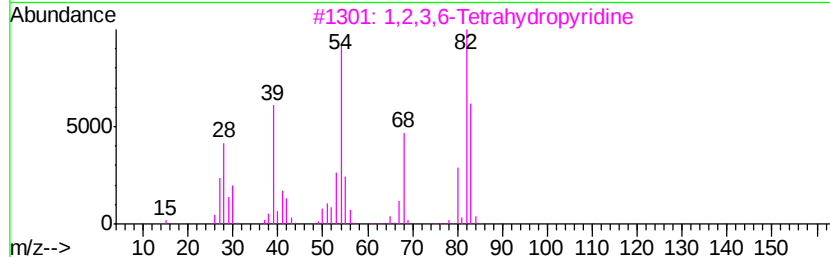
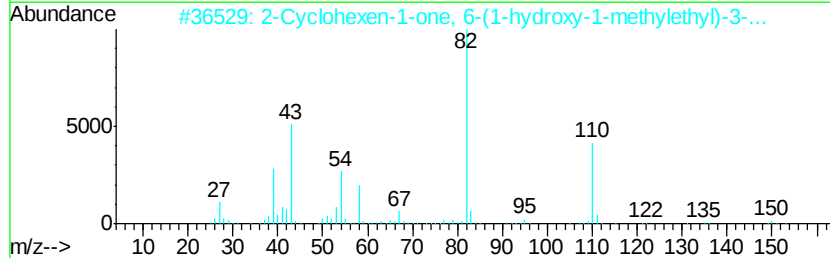
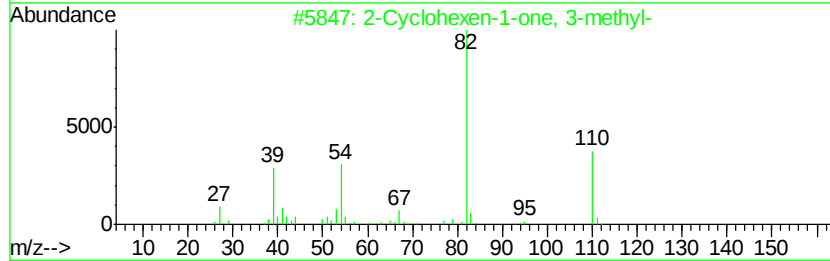
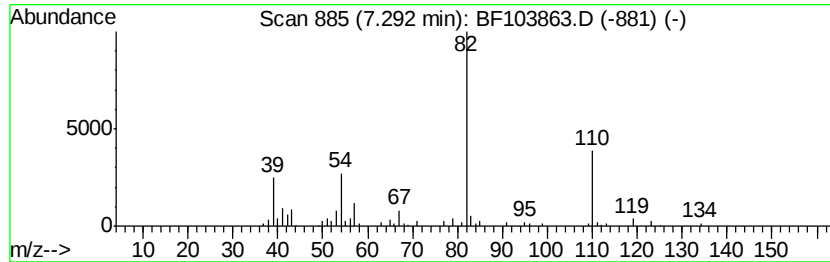
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Cyclohexen-1-one, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.29	2.71 ng	116221	1,4-Dichlorobenzene-d4	6.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-one, 3-methyl-	110	C7H10O	001193-18-6	81
2	2-Cyclohexen-1-one, 6-(1-hydroxy...	168	C10H16O2	087791-00-2	56
3	1,2,3,6-Tetrahydropyridine	83	C5H9N	000694-05-3	47
4	2-Cyclohexen-1-one, 3,5-dimethyl-	124	C8H12O	001123-09-7	43
5	1H-Pyrazole, 3-methyl-	82	C4H6N2	001453-58-3	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
Data File : BF103863.D
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Operator : JU/SJ
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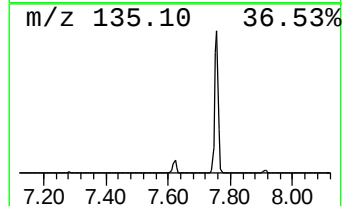
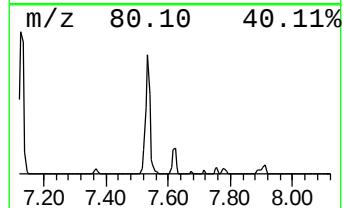
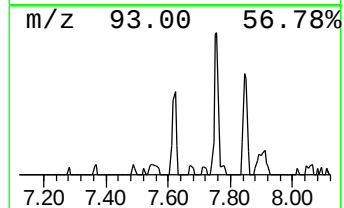
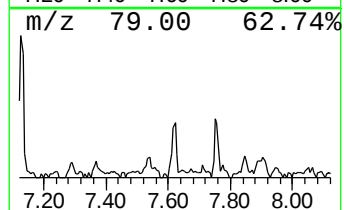
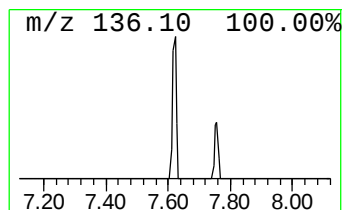
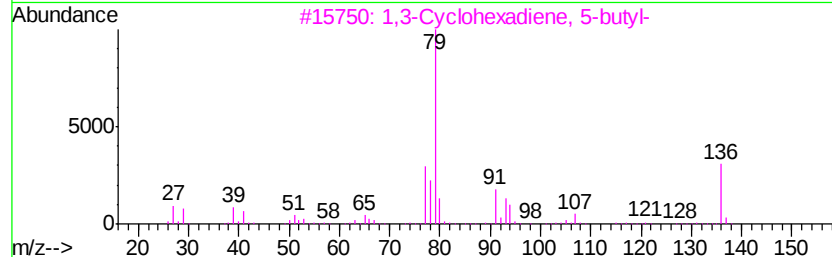
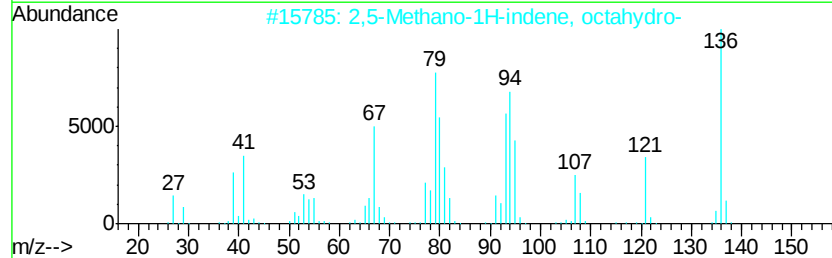
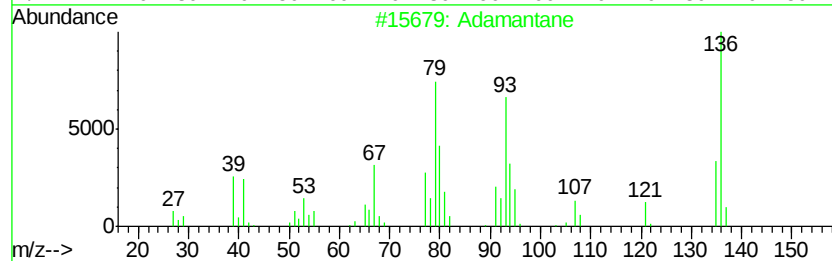
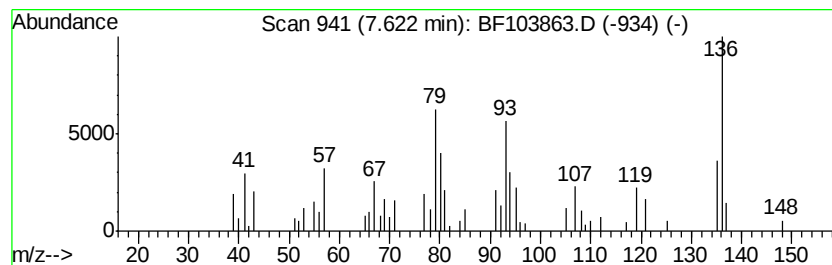
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Adamantane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	3.07 ng	131455	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Adamantane	136	C10H16	000281-23-2	95
2		2,5-Methano-1H-indene, octahydro-	136	C10H16	019026-94-9	46
3		1,3-Cyclohexadiene, 5-butyl-	136	C10H16	030168-57-1	46
4		Tricyclo[3.3.2.0(3,7)]decane	136	C10H16	049700-60-9	43
5		1H-Inden-1-one, 2,3,4,5,6,7-hexa...	136	C9H12O	022118-00-9	38



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF032218\
Data File : BF103863.D
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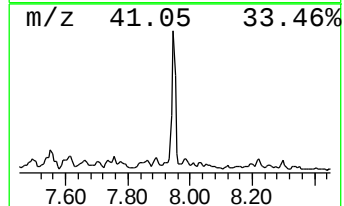
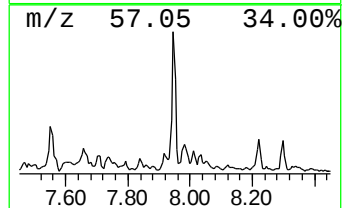
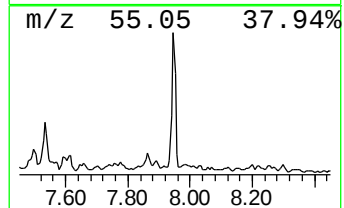
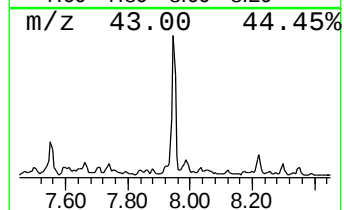
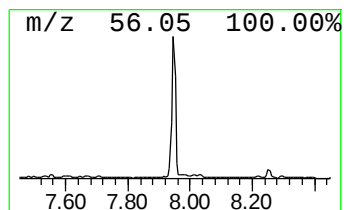
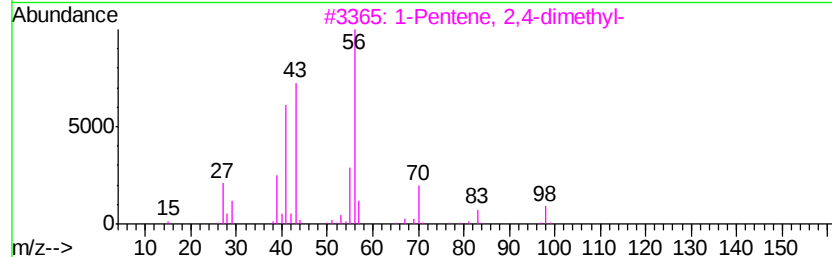
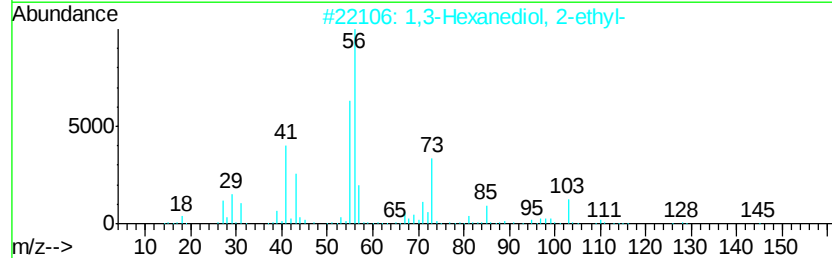
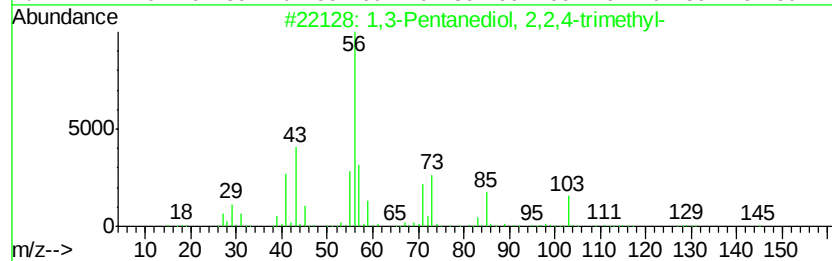
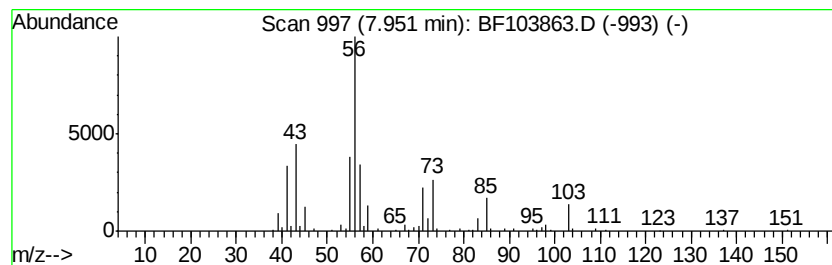
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,3-Pentanediol, 2,2,4-trim... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	6.49 ng	347623	Napthalene-d8	8.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	90
2			1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	40
3			1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	38
4			1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38
5			Cyclopentanamine	85	C5H11N	001003-03-8	37



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Data File : BF103863.D
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Operator : JU/SJ
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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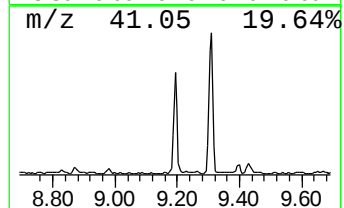
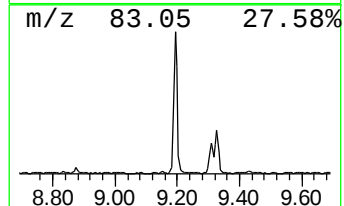
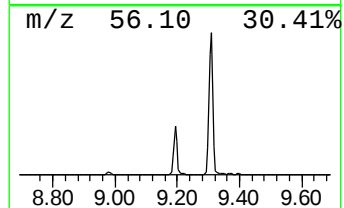
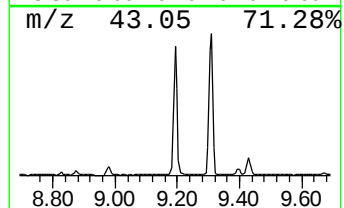
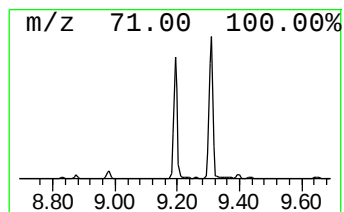
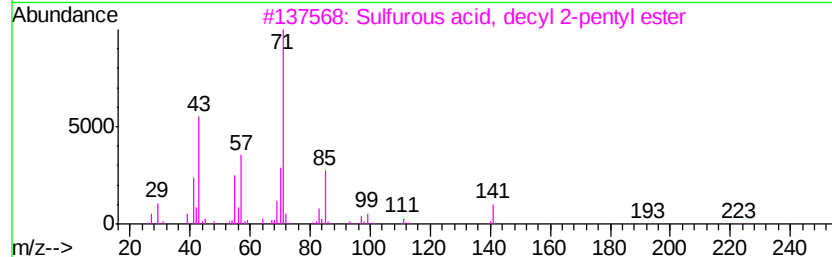
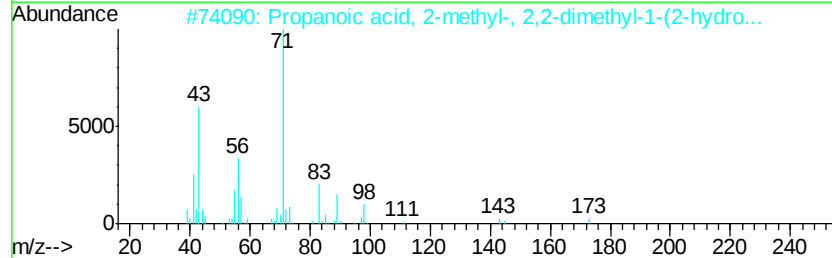
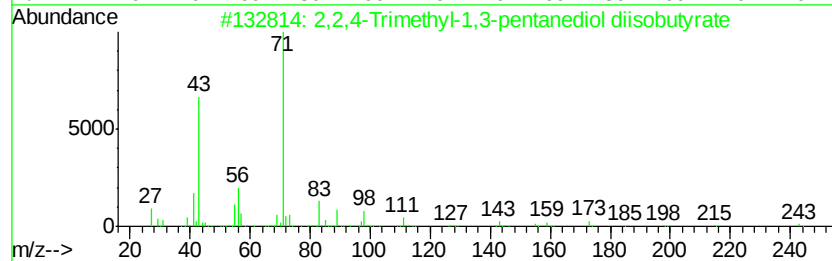
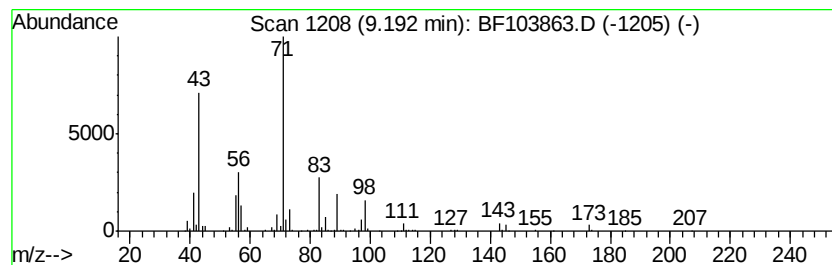
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.19	10.52 ng	548036	Acenaphthene-d10	10.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	59
2		Propanoic acid, 2-methyl-, 2,2-d...	216	C12H24O3	074367-33-2	56
3		Sulfurous acid, decyl 2-pentyl e...	292	C15H32O3S	1000309-15-9	43
4		Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	42
5		Butanoic acid, 1-methyloctyl ester	214	C13H26O2	069727-42-0	40



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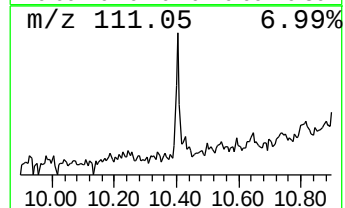
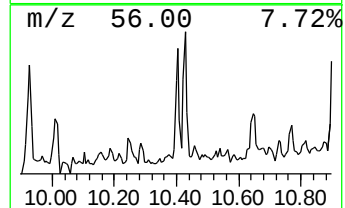
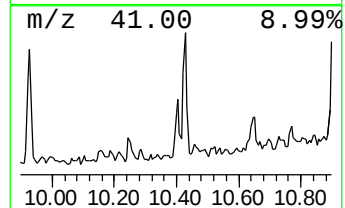
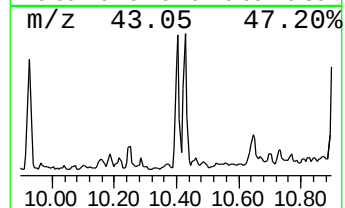
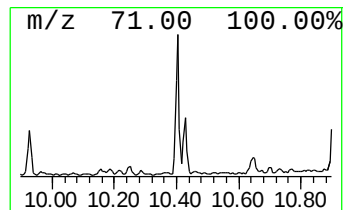
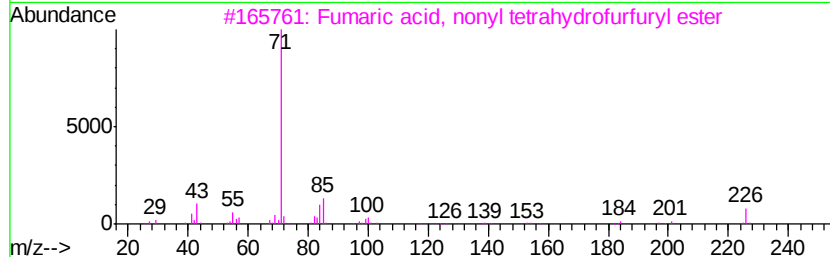
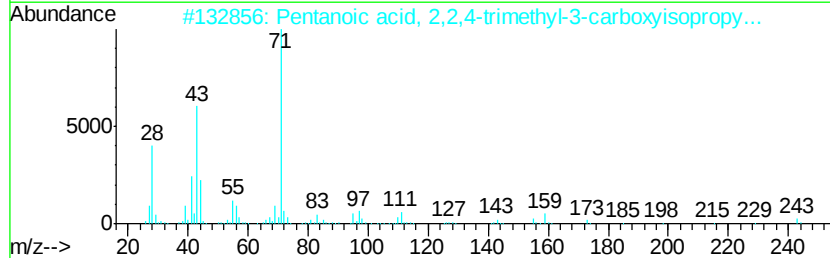
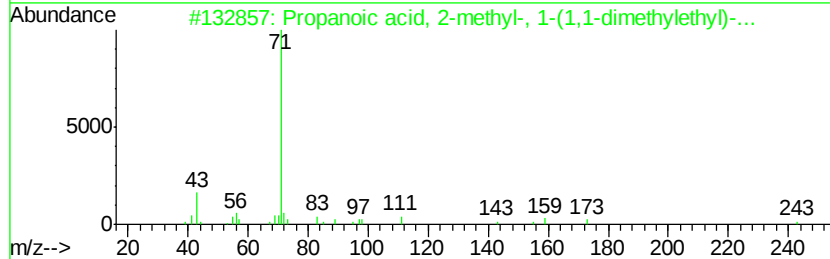
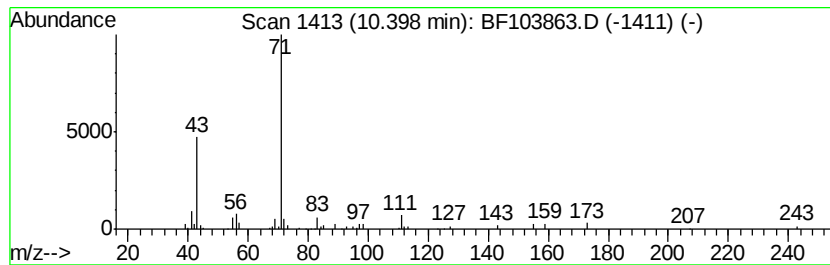
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Propanoic acid, 2-methyl-, ... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.40	2.01 ng	104526	Acenaphthene-d10	10.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Propanoic acid, 2-methyl-, 1-(1,...	286 C16H3004	074381-40-1 90
2		Pentanoic acid, 2,2,4-trimethyl-...	286 C16H3004	1000140-77-5 74
3		Fumaric acid, nonyl tetrahydrofu...	326 C18H3005	1000330-58-3 59
4		4-Hexen-3-ol, 2-methyl-	114 C7H140	004798-60-1 59
5		2,2,4-Trimethyl-1,3-pentanediol ...	286 C16H3004	006846-50-0 56



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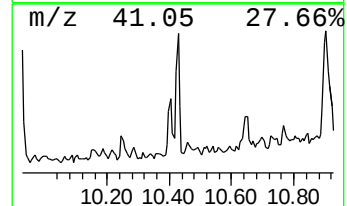
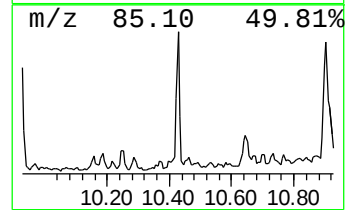
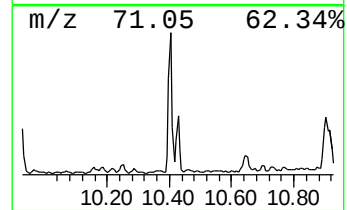
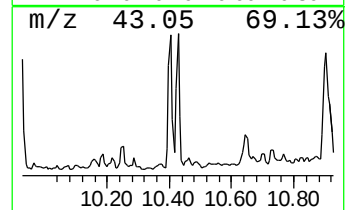
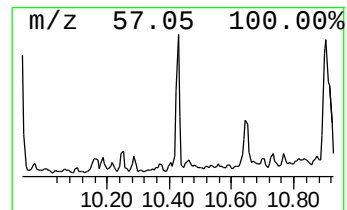
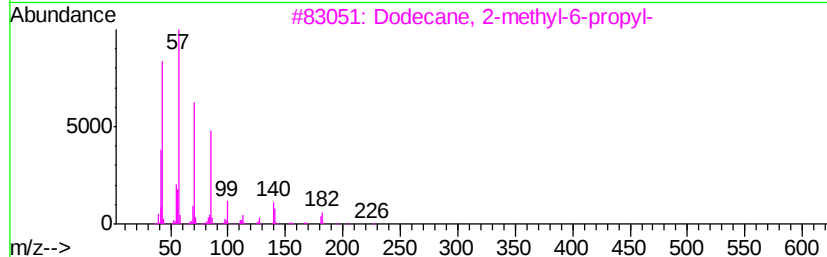
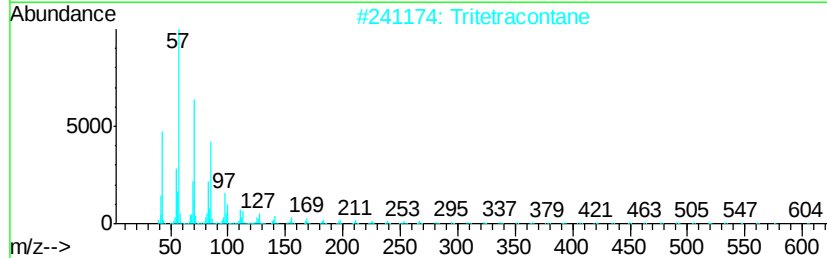
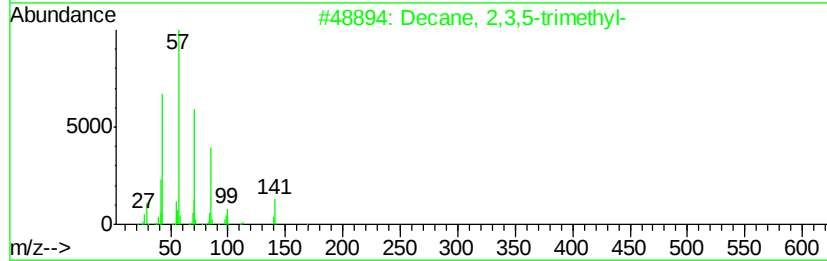
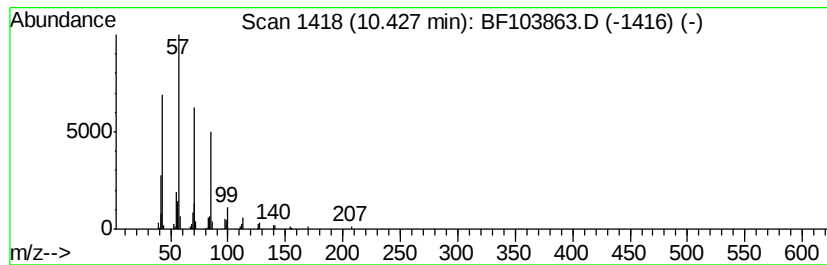
Instrument :
BNA_F
ClientSampleId :
SLAB-CONCRETE-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Decane, 2,3,5-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	2.11 ng	109794	Acenaphthene-d10	10.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3	90
2	Tritetracontane	605 C43H88	007098-21-7	86
3	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	86
4	Dodecane, 1-iodo-	296 C12H25I	004292-19-7	86
5	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	78



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
Data File : BF103863.D
Acq On : 22 Mar 2018 20:18
Operator : JU/SJ
Sample : J1997-25
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SLAB-CONCRETE-B

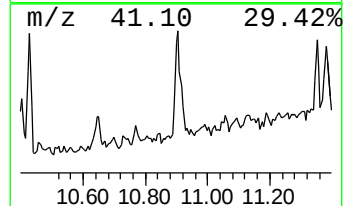
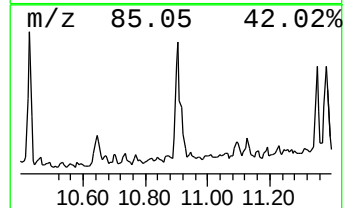
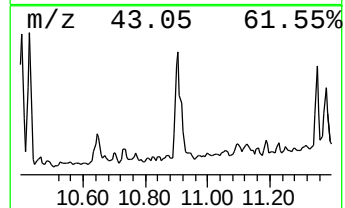
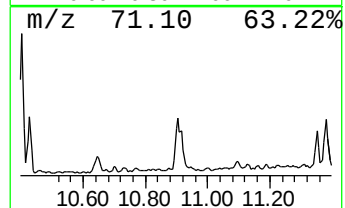
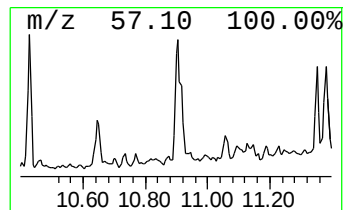
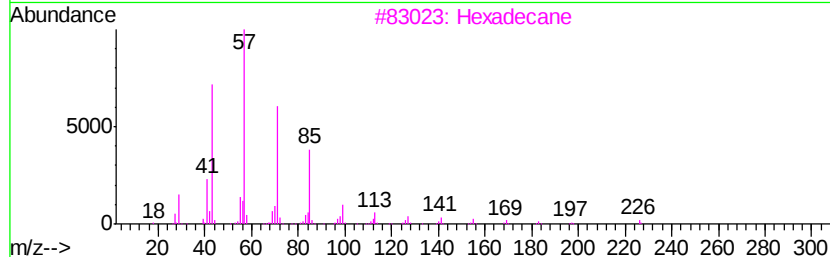
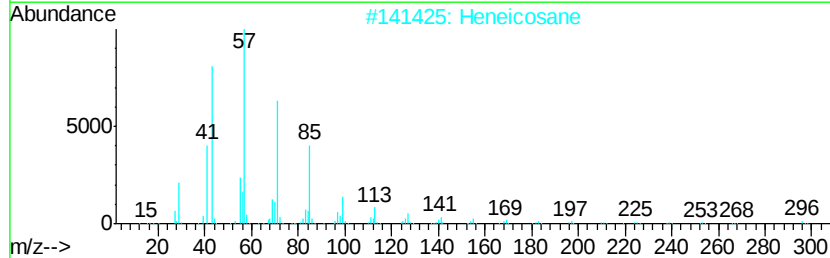
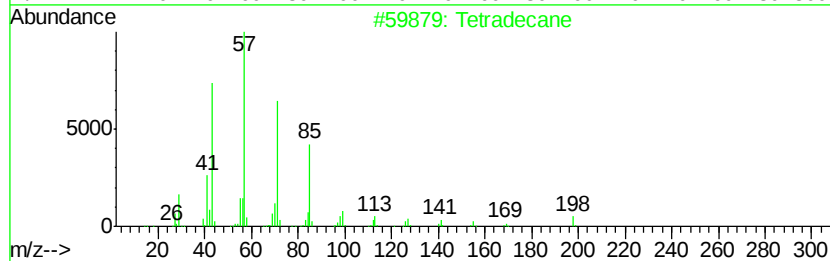
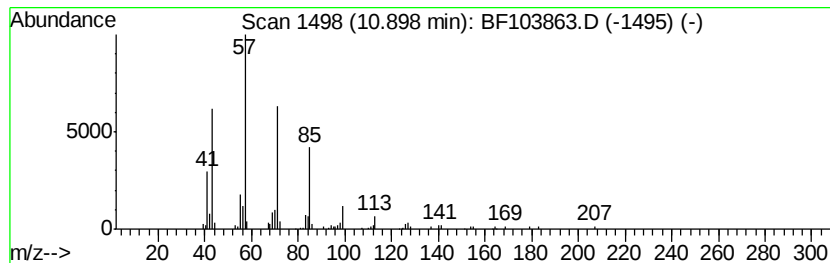
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	4.18 ng	178168	Phenanthrene-d10	11.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Hexadecane	226	C16H34	000544-76-3	91
4			Pentadecane	212	C15H32	000629-62-9	91
5			Tritetracontane	605	C43H88	007098-21-7	90



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
Data File : BF103863.D
Acq On : 22 Mar 2018 20:18
Operator : JU/SJ
Sample : J1997-25
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SLAB-CONCRETE-B

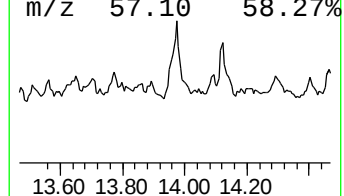
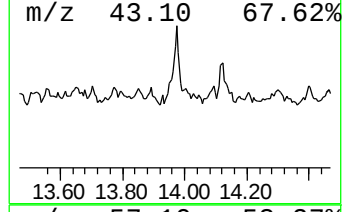
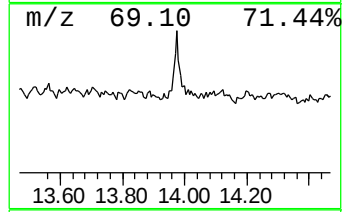
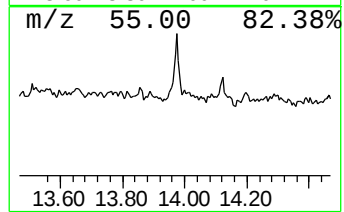
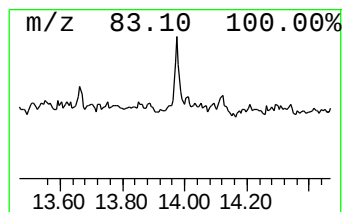
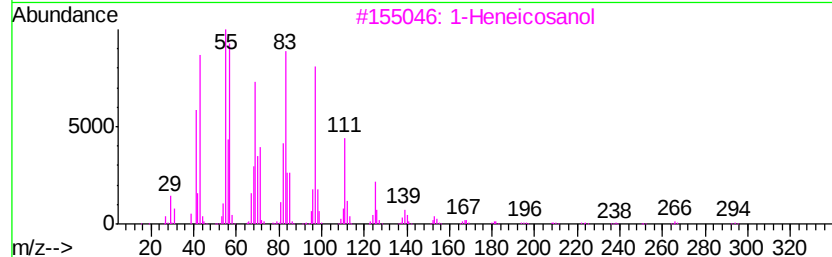
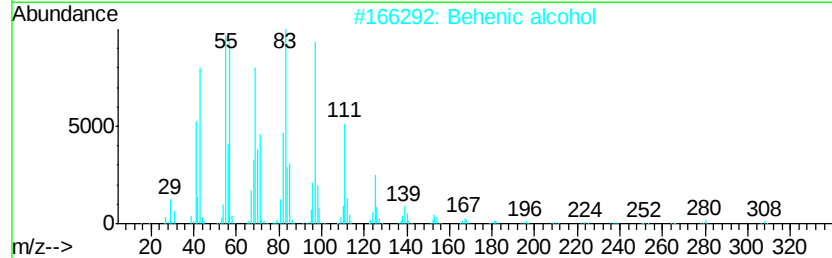
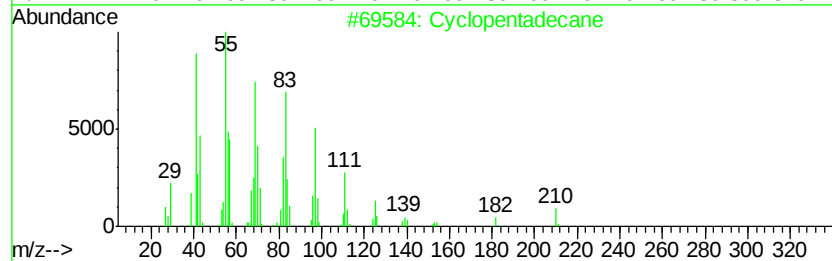
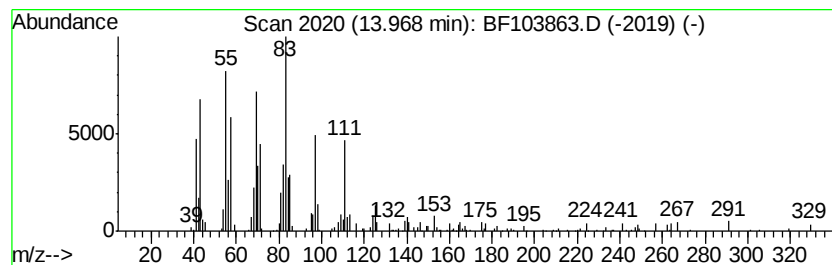
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclopentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	4.58 ng	186219	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	94
2		Behenic alcohol	326	C22H46O	000661-19-8	93
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	90
5		1-Octadecene	252	C18H36	000112-88-9	90



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
 Data File : BF103863.D
 Acq On : 22 Mar 2018 20:18
 Operator : JU/SJ
 Sample : J1997-25
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SLAB-CONCRETE-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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...	2.30	3.4	ng	146714	1	6.97	857684	20.0
4-Penten-2-one, 4...	3.79	2.8	ng	121731	1	6.97	857684	20.0
3-Penten-2-one, 4...	4.63	119.9	ng	5140530	1	6.97	857684	20.0
2-Pentanone, 4-hy...	5.24	154.0	ng	6606060	1	6.97	857684	20.0
unknown6.74	6.74	17.3	ng	741675	1	6.97	857684	20.0
3-Heptene, 2,2,4,...	6.87	2.0	ng	86822	1	6.97	857684	20.0
2-Cyclohexen-1-on...	7.29	2.7	ng	116221	1	6.97	857684	20.0
Adamantane	7.62	3.1	ng	131455	1	6.97	857684	20.0
1,3-Pentanediol, ...	7.95	6.5	ng	347623	2	8.26	1071270	20.0
2,2,4-Trimethyl-1...	9.19	10.5	ng	548036	3	10.02	1041930	20.0
Propanoic acid, 2...	10.40	2.0	ng	104526	3	10.02	1041930	20.0
Decane, 2,3,5-tri...	10.43	2.1	ng	109794	3	10.02	1041930	20.0
Tetradecane	10.90	4.2	ng	178168	4	11.50	852283	20.0
Cyclopentadecane	13.97	4.6	ng	186219	5	14.16	813470	20.0