

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
 Data File : BF077505.D
 Acq On : 27 Feb 2015 14:35
 Operator : TP/IZ
 Sample : G1377-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RW-3

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: CENTROID

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.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	1.550	54	58	61	rVB	365210	487472	20.80%	2.676%
2	1.721	70	73	76	rBV	198677	273660	11.68%	1.502%
3	1.847	82	84	92	rVB	158033	224132	9.57%	1.231%
4	5.036	361	363	367	rVB	93424	105968	4.52%	0.582%
5	5.550	405	408	410	rBV	524352	624202	26.64%	3.427%
6	6.807	516	518	521	rBV	303105	350031	14.94%	1.922%
7	6.967	529	532	534	rBV	1269281	1136196	48.49%	6.238%
8	7.253	555	557	560	rBV	374824	443886	18.94%	2.437%
9	7.447	571	574	575	rBV	1121204	1122235	47.90%	6.161%
10	7.470	575	576	578	rVB	50170	68138	2.91%	0.374%
11	7.950	613	618	620	rBV	970181	1009064	43.07%	5.540%
12	8.213	636	641	643	rBV	59760	64221	2.74%	0.353%
13	8.430	656	660	665	rBV2	30808	73632	3.14%	0.404%
14	8.522	665	668	670	rVB2	155095	191988	8.19%	1.054%
15	8.830	693	695	697	rVB	617965	675670	28.84%	3.710%
16	8.865	697	698	700	rVV2	45400	55555	2.37%	0.305%
17	8.899	700	701	704	rVB	71311	67550	2.88%	0.371%
18	9.185	724	726	728	rVB2	23483	31097	1.33%	0.171%
19	9.322	735	738	740	rVB	39057	43973	1.88%	0.241%
20	9.425	745	747	749	rVV	48738	56499	2.41%	0.310%
21	9.470	749	751	754	rVV3	16696	35869	1.53%	0.197%
22	9.539	755	757	758	rVV2	51015	47192	2.01%	0.259%
23	9.573	758	760	762	rVV	28348	27590	1.18%	0.151%
24	9.653	766	767	769	rVB	39709	49433	2.11%	0.271%
25	9.722	769	773	775	rBV	68846	79999	3.41%	0.439%
26	9.836	781	783	785	rBV	275285	296574	12.66%	1.628%
27	9.893	785	788	791	rVB2	24859	42072	1.80%	0.231%
28	9.985	791	796	797	rBV2	14270	32925	1.41%	0.181%
29	10.179	811	813	816	rVB	2090456	1876376	80.08%	10.302%
30	10.351	825	828	832	rBV3	13307	45006	1.92%	0.247%
31	10.431	832	835	837	rVB	31701	44113	1.88%	0.242%
32	10.499	838	841	843	rBV	67182	101010	4.31%	0.555%
33	10.591	846	849	851	rBV	127882	156608	6.68%	0.860%
34	10.625	851	852	855	rVB2	73686	63241	2.70%	0.347%

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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: CENTROID

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	10.728	860	861	865	rVB2	40271	64325	2.75%	0.353%
36	10.831	869	870	872	rBV	36012	44338	1.89%	0.243%
37	10.922	874	878	880	rVB2	12165	25761	1.10%	0.141%
38	11.002	883	885	888	rBV	591313	826595	35.28%	4.538%
39	11.151	894	898	901	rBV4	14359	44135	1.88%	0.242%
40	11.265	905	908	910	rVB2	28733	34934	1.49%	0.192%
41	11.322	910	913	915	rBV3	17690	27963	1.19%	0.154%
42	11.414	918	921	922	rBV	33639	45105	1.93%	0.248%
43	11.539	925	932	933	rBV	46930	84099	3.59%	0.462%
44	11.562	933	934	937	rVB	60633	47710	2.04%	0.262%
45	11.631	937	940	943	rVB5	10536	28129	1.20%	0.154%
46	11.688	943	945	947	rBV	48570	50224	2.14%	0.276%
47	11.768	950	952	953	rBV	39143	33667	1.44%	0.185%
48	11.905	961	964	967	rBV	85188	118764	5.07%	0.652%
49	11.985	969	971	974	rVB	1238383	1187048	50.66%	6.517%
50	12.145	981	985	991	rVB7	10720	29328	1.25%	0.161%
51	12.374	1000	1005	1006	rBV3	17432	32315	1.38%	0.177%
52	12.442	1009	1011	1013	rVB2	24055	30851	1.32%	0.169%
53	12.694	1028	1033	1035	rBV5	16433	33186	1.42%	0.182%
54	12.854	1044	1047	1049	rBV	777437	876952	37.43%	4.815%
55	13.562	1107	1109	1111	rBV3	27853	35866	1.53%	0.197%
56	14.842	1219	1221	1224	rVB	1937201	2343107	100.00%	12.864%
57	16.134	1332	1334	1337	rBV	887790	1067686	45.57%	5.862%
58	17.231	1427	1430	1435	rBV6	13401	46802	2.00%	0.257%
59	17.883	1484	1487	1490	rBV	930612	1051796	44.89%	5.775%

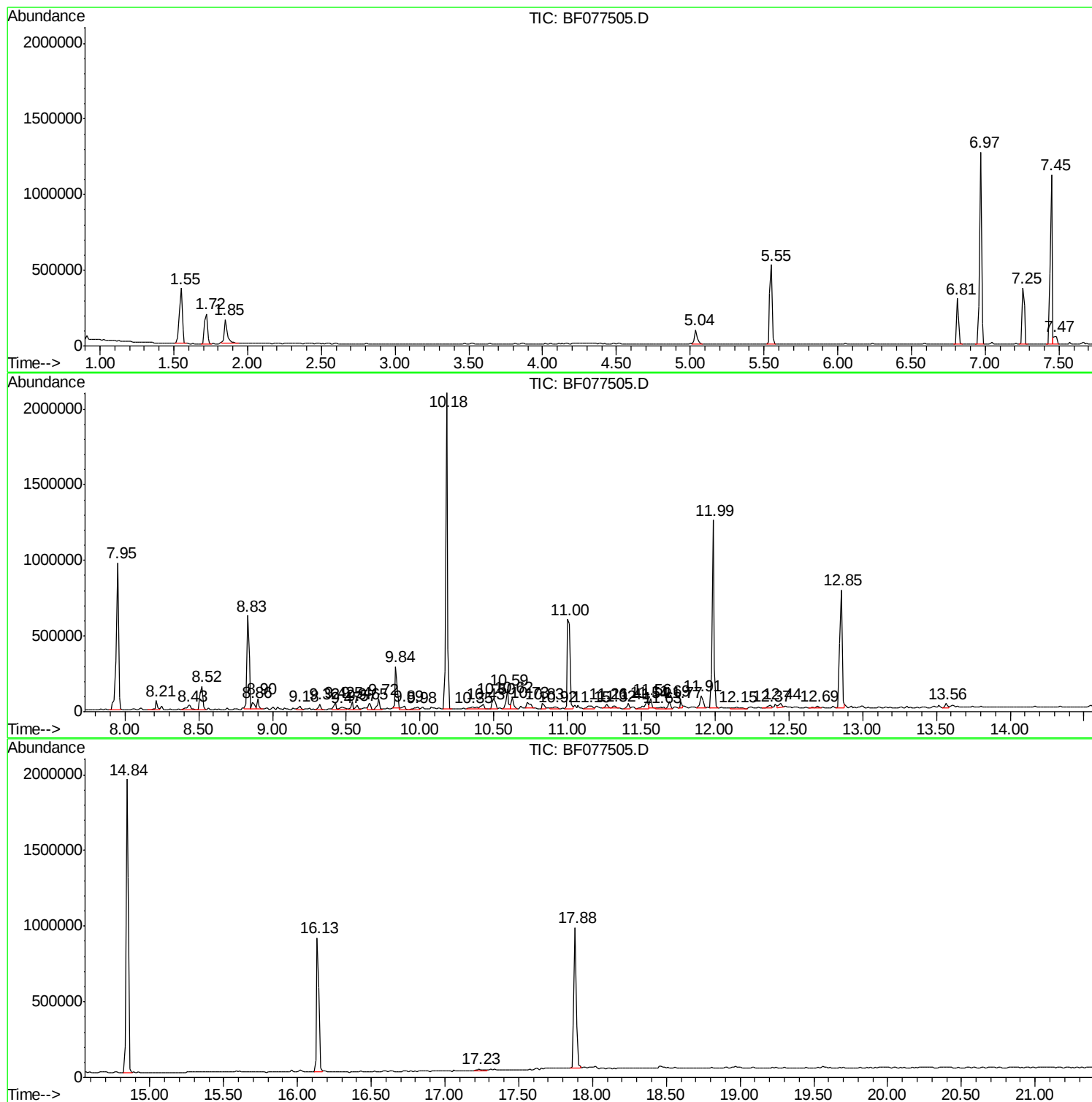
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ClientSampled :
RW-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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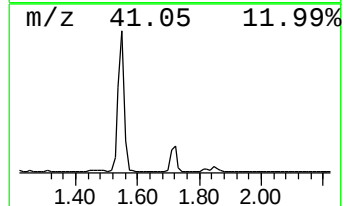
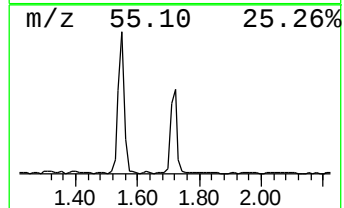
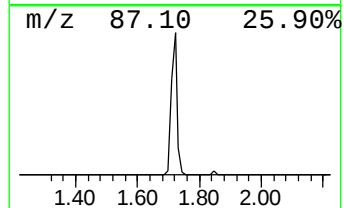
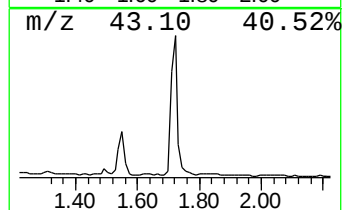
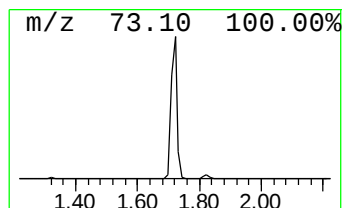
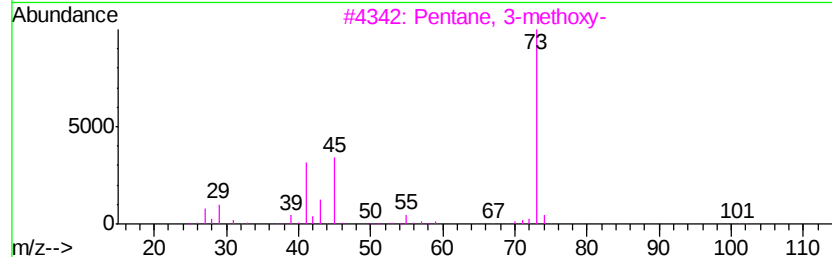
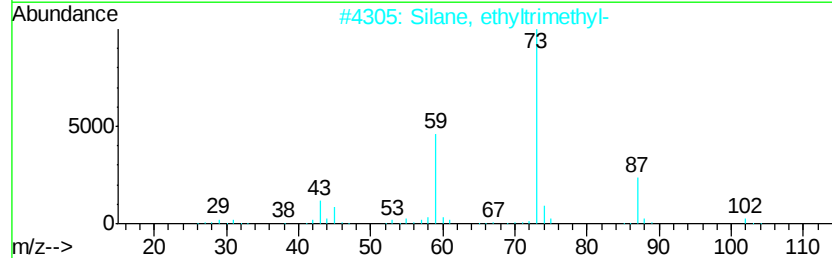
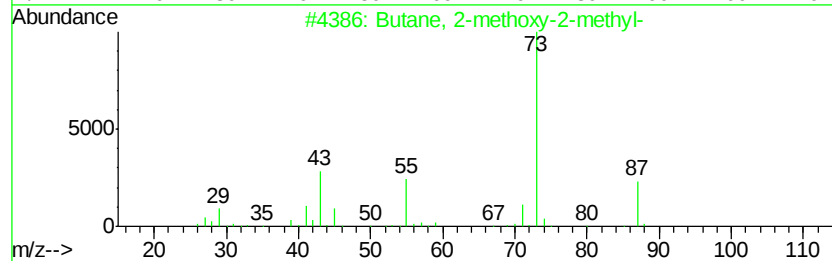
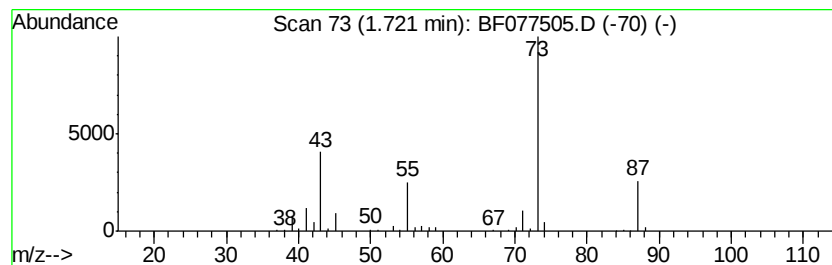
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.72	12.33 ng	273660	1,4-Dichlorobenzene-d4	7.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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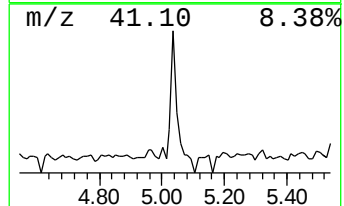
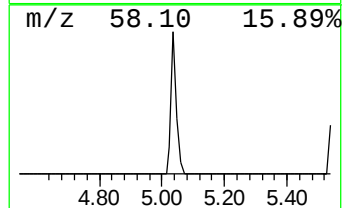
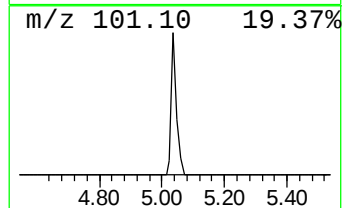
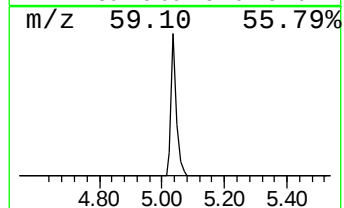
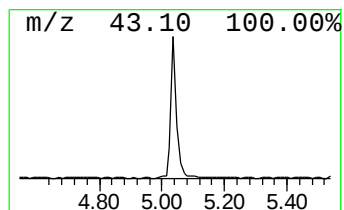
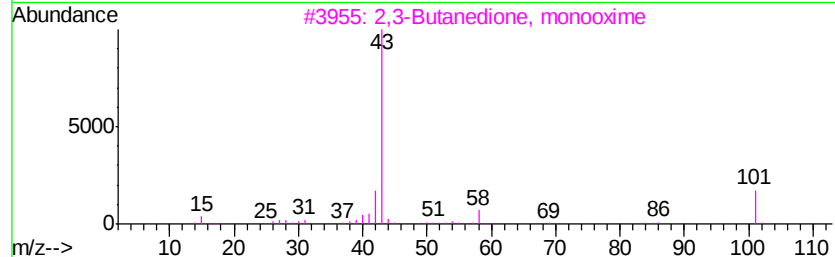
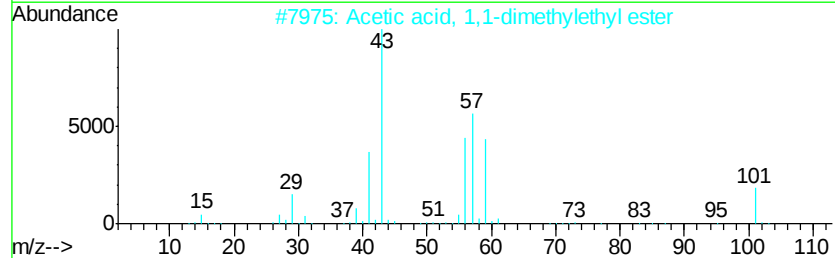
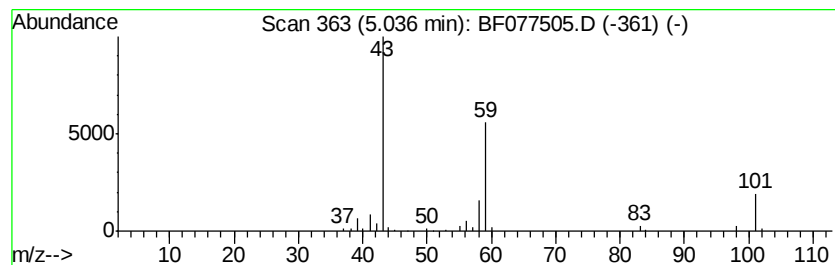
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	4.77 ng	105968	1,4-Dichlorobenzene-d4	7.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		5-Hexen-2-one	98	C6H10O	000109-49-9	10



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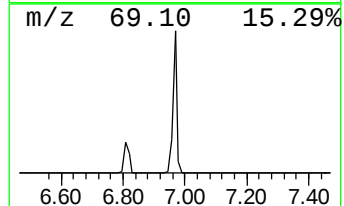
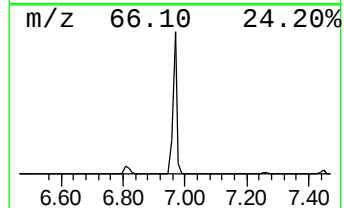
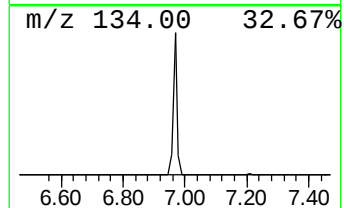
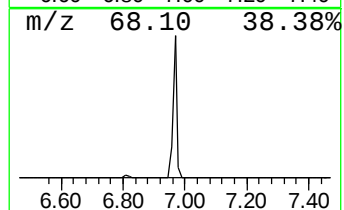
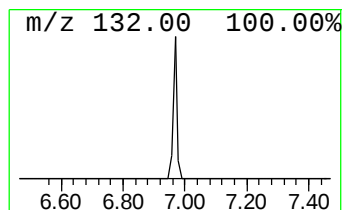
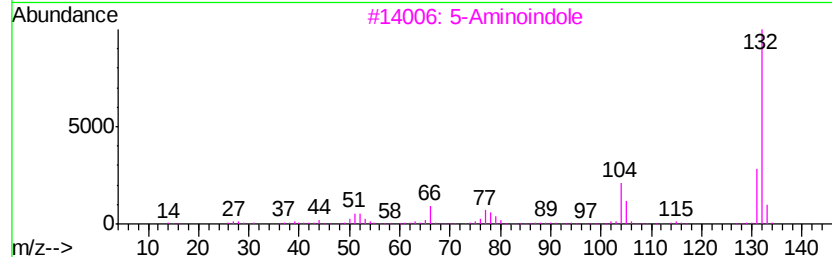
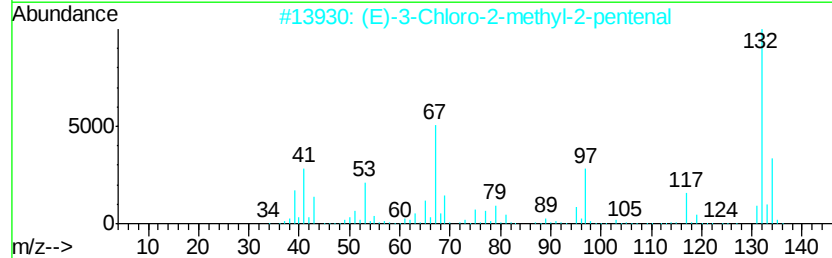
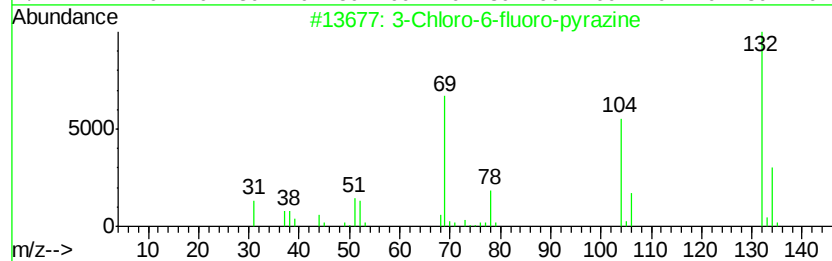
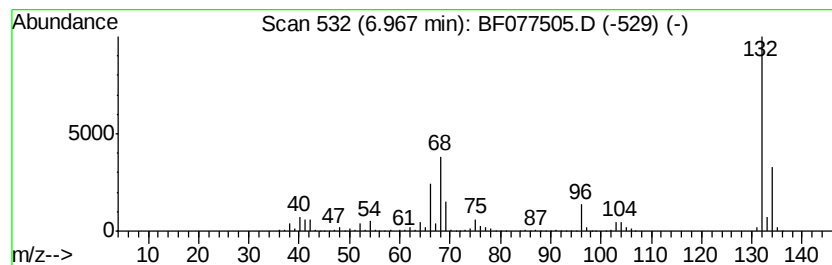
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.97 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.97	51.19 ng	1136200	1,4-Dichlorobenzene-d4	7.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	16
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
4	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9
5	4-Methylpyrrolo[1,2-a]pyrazine			132	C8H8N2	064608-60-2	9



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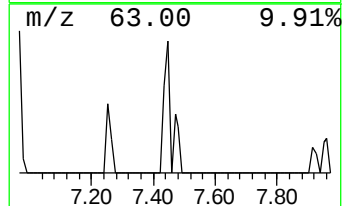
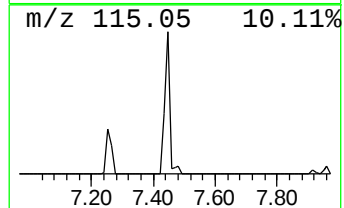
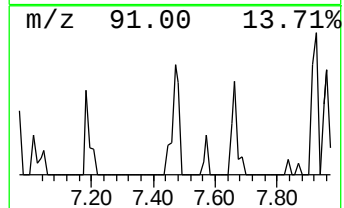
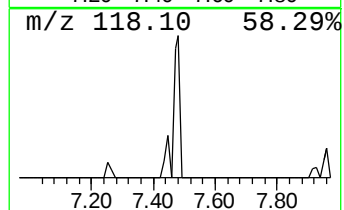
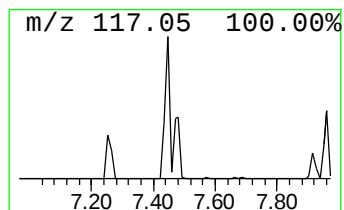
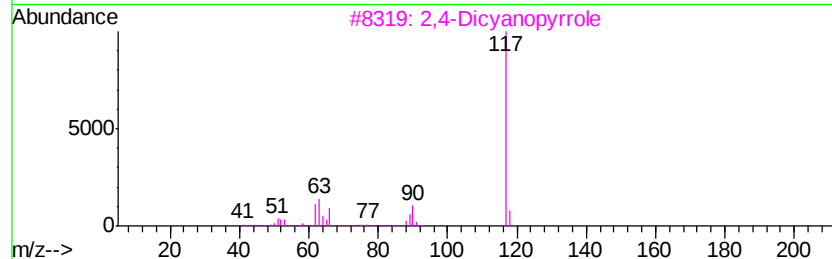
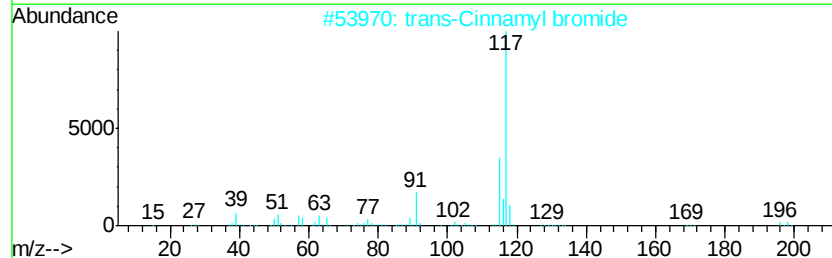
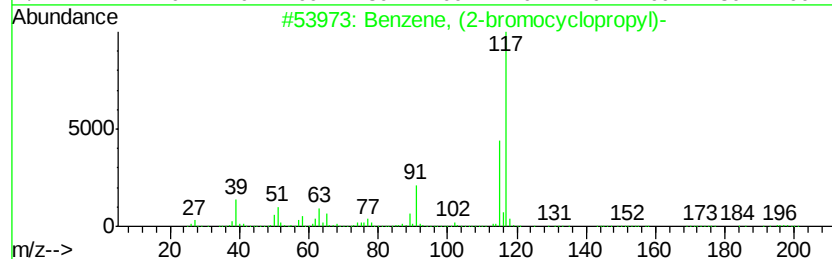
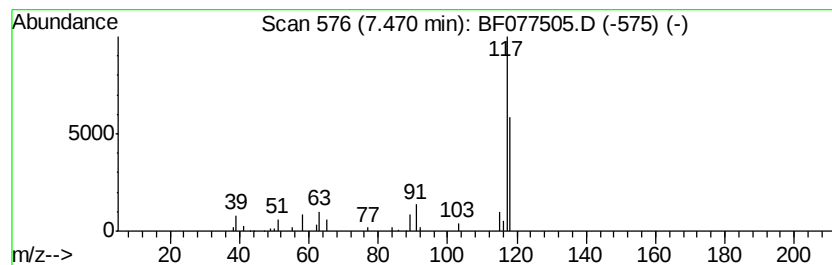
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown7.47 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	3.07 ng	68138	1,4-Dichlorobenzene-d4	7.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	28
2		trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	28
3		2,4-Dicyanopyrrole	117	C6H3N3	074023-87-3	25
4		Indole	117	C8H7N	000120-72-9	9
5		1H-Pyrrole-3,4-dicarbonitrile	117	C6H3N3	125666-25-3	9



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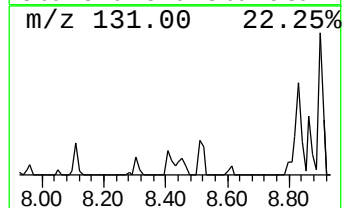
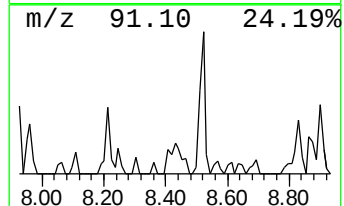
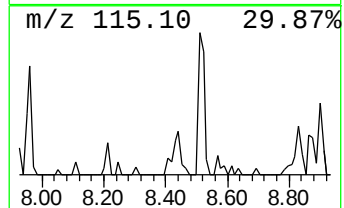
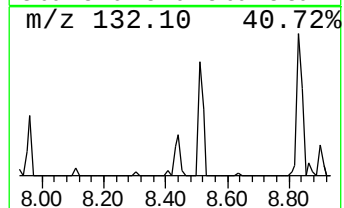
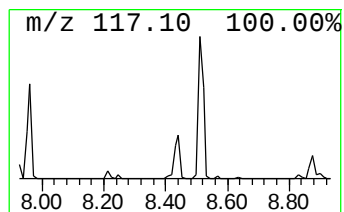
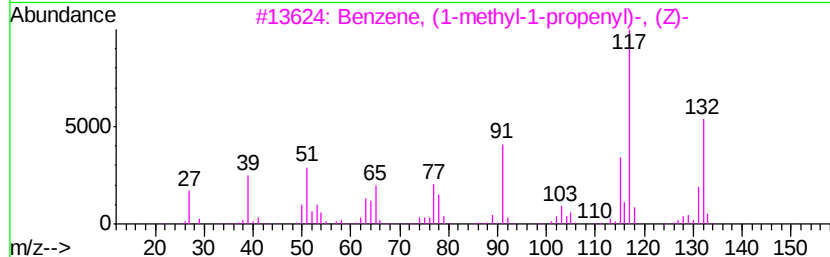
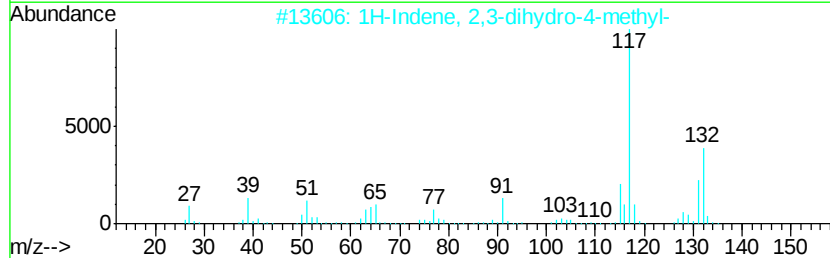
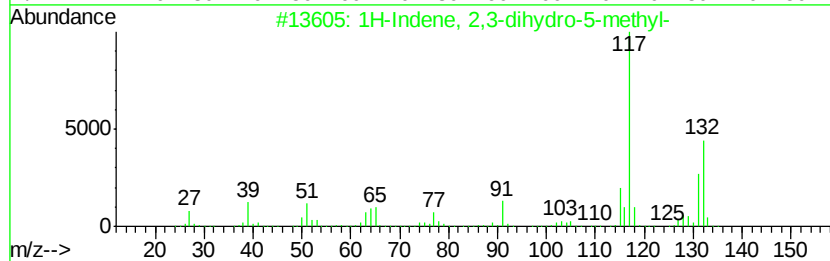
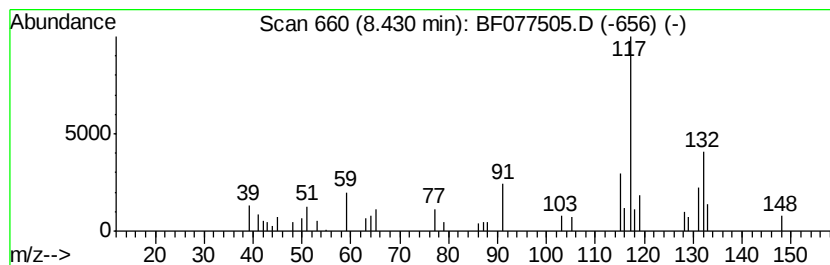
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.43	2.18 ng	73632	Napthalene-d8	8.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	68
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
Operator : TP/IZ
Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-3

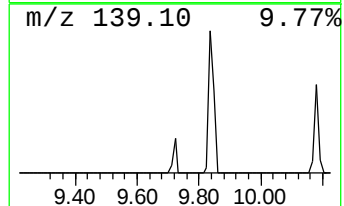
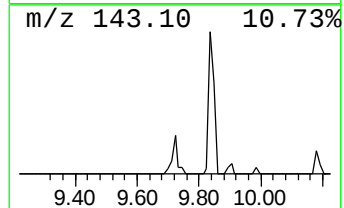
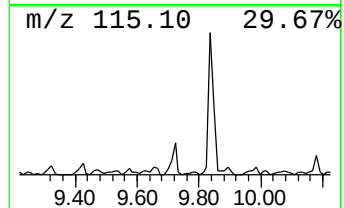
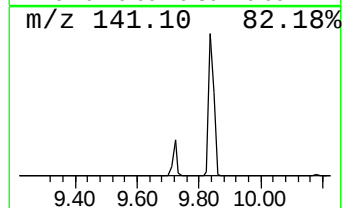
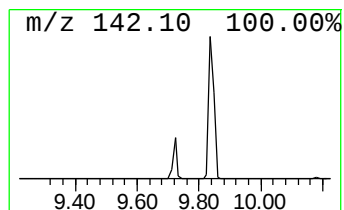
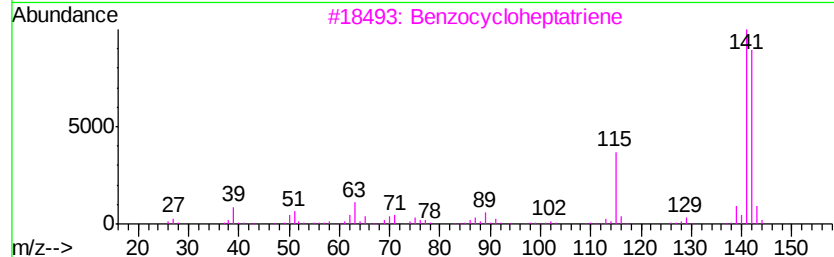
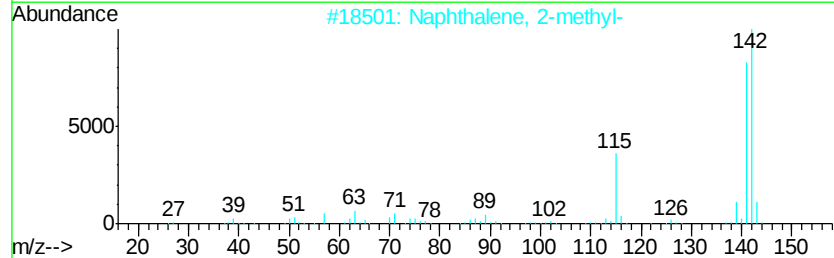
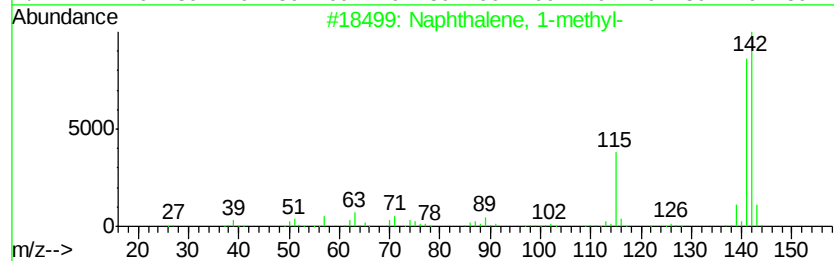
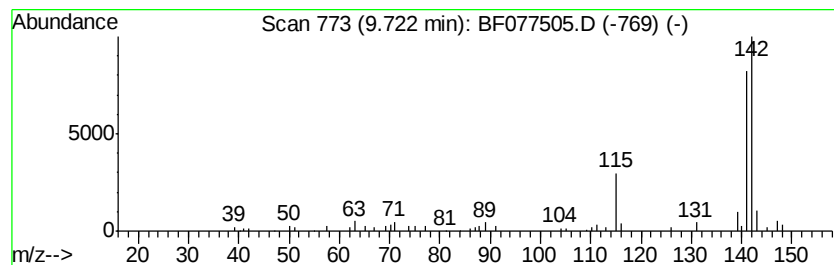
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	2.37 ng	79999	Naphthalene-d8	8.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
3		Benzocycloheptatriene	142	C11H10	000264-09-5	86
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	86
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
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Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RW-3

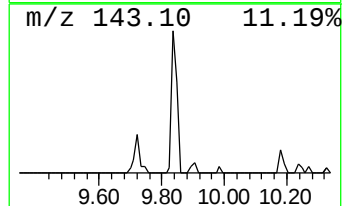
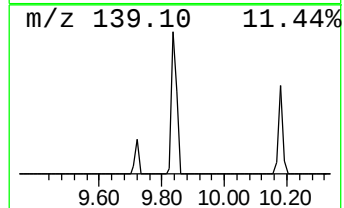
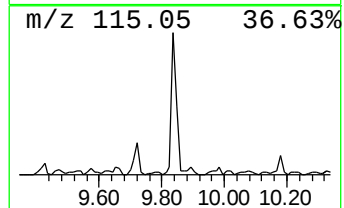
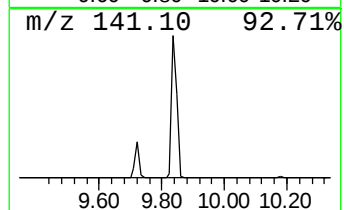
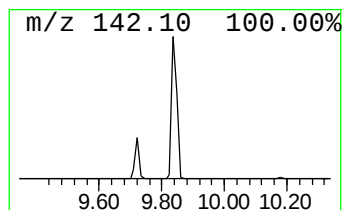
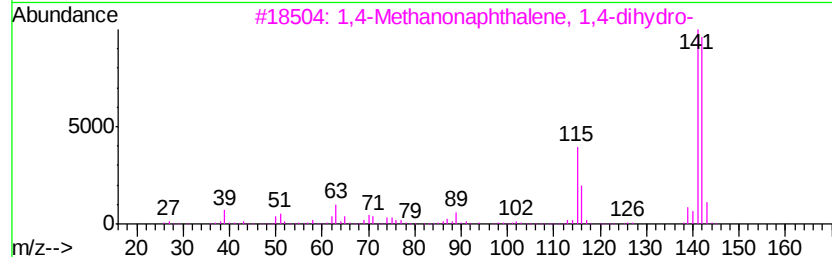
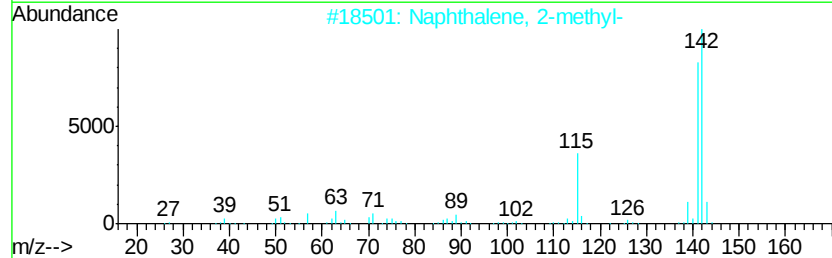
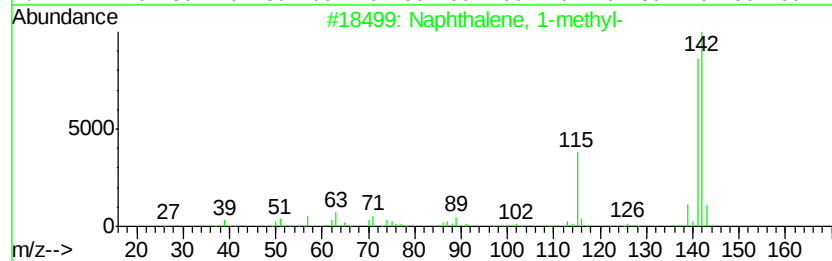
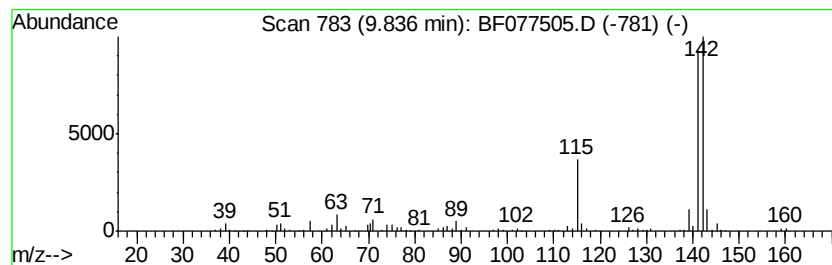
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Naphthalene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.84	8.78 ng	296574	Naphthalene-d8	8.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
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Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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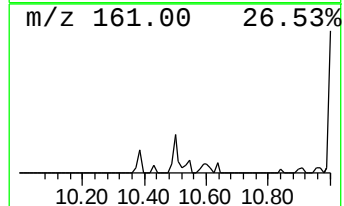
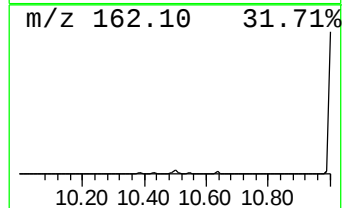
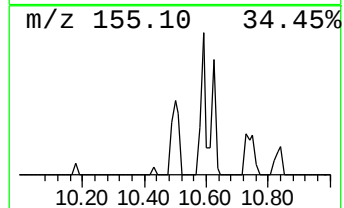
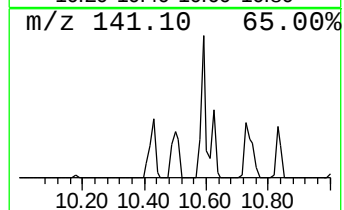
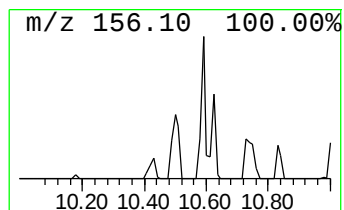
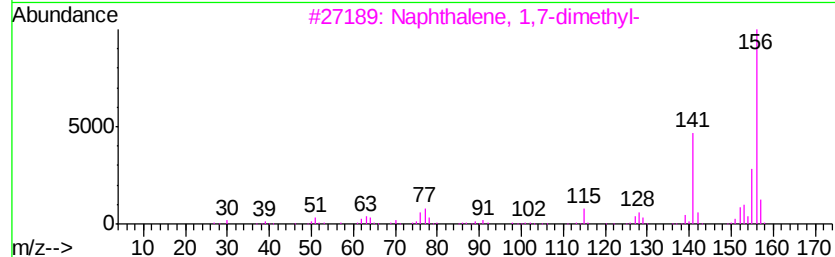
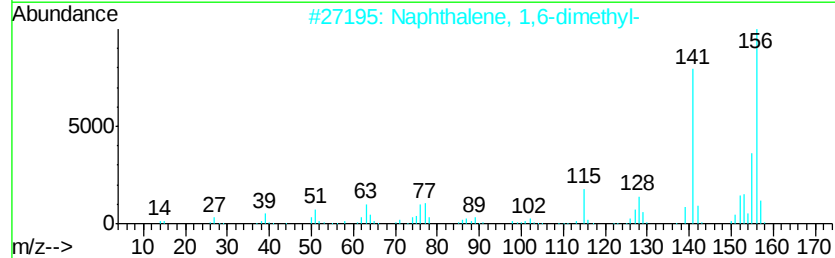
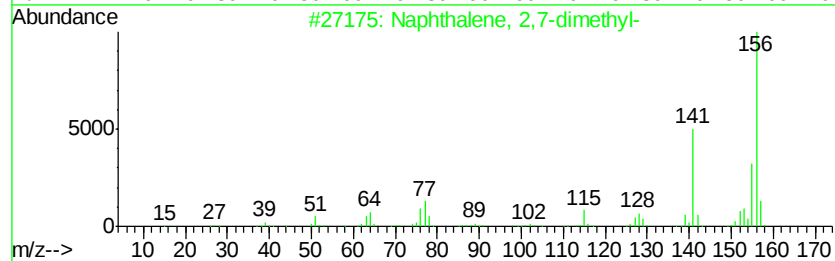
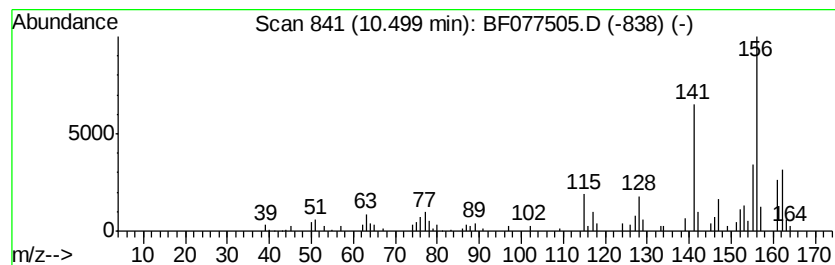
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Naphthalene, 2,7-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.50	2.44 ng	101010	Acenaphthene-d10	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
Operator : TP/IZ
Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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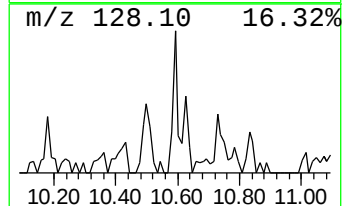
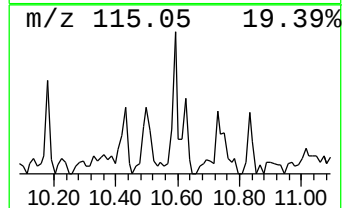
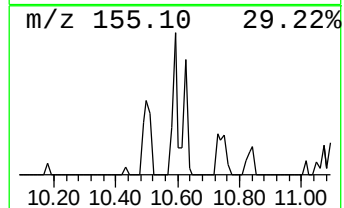
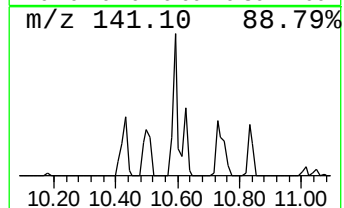
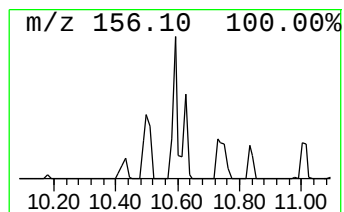
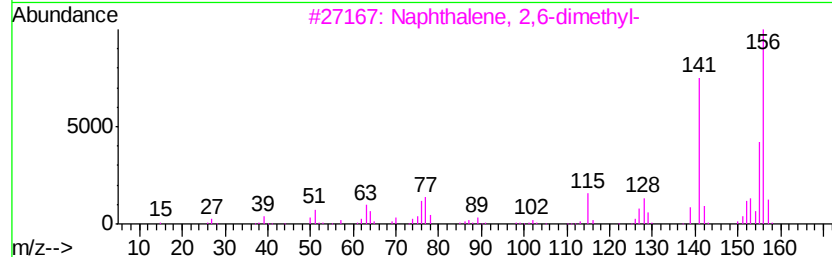
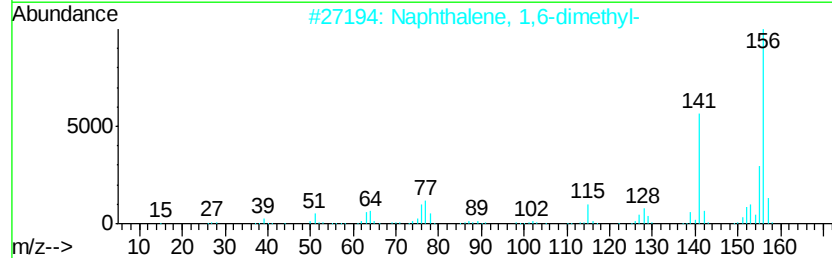
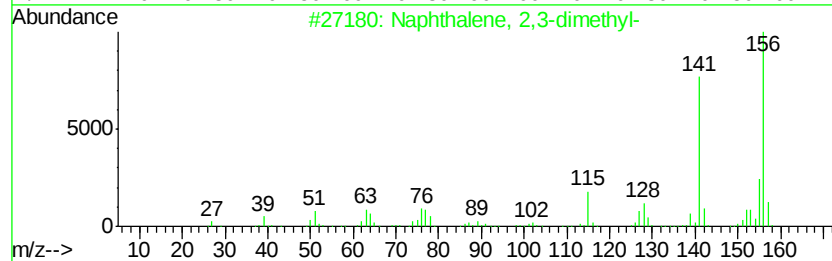
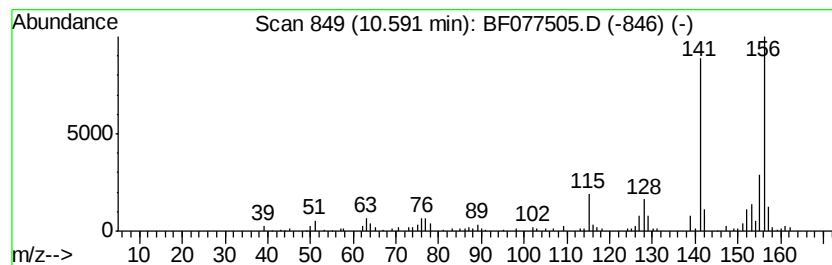
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.59	3.79 ng	156608	Acenaphthene-d10	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
Operator : TP/IZ
Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-3

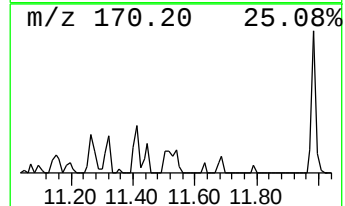
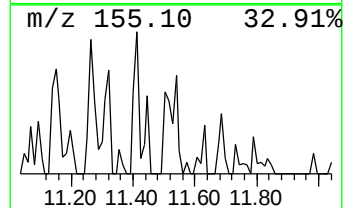
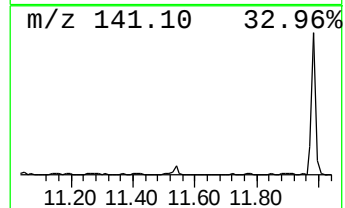
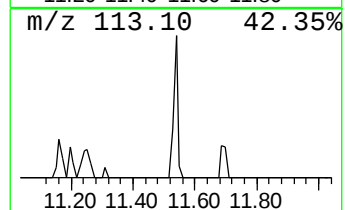
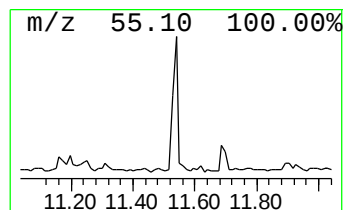
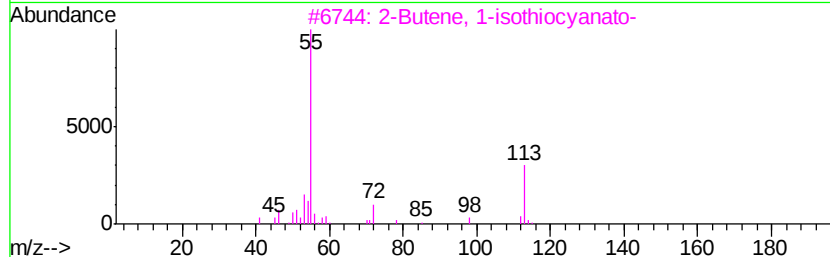
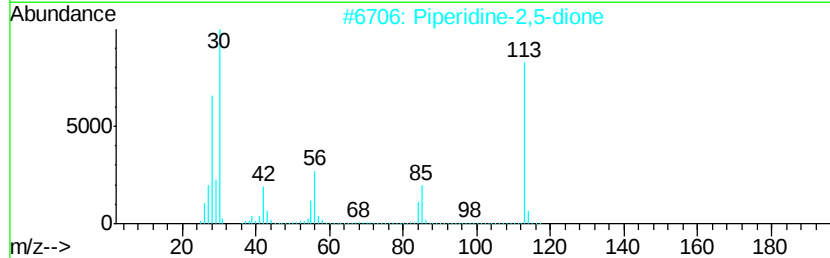
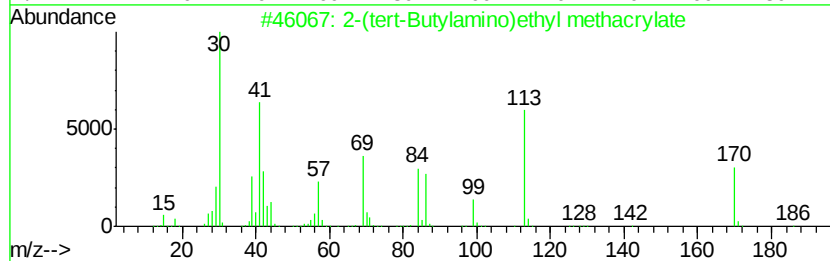
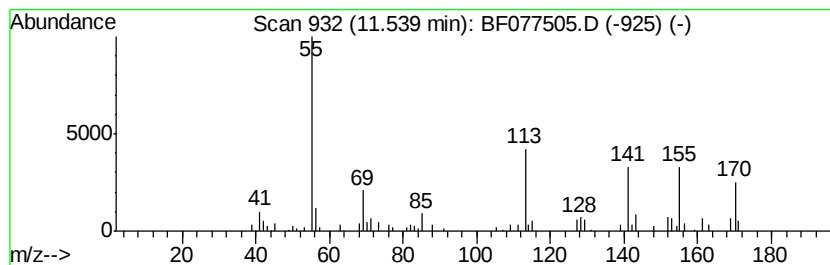
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown11.54 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.54	2.03 ng	84099	Acenaphthene-d10	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(tert-Butylamino)ethyl methacr...	185	C10H19N02	003775-90-4	17
2		Piperidine-2,5-dione	113	C5H7N02	052065-78-8	10
3		2-Butene, 1-isothiocyanato-	113	C5H7NS	002253-93-2	9
4		3,4-Dimethyl-5-hydroxy-isoxazole	113	C5H7N02	029279-99-0	9
5		3,4-Dimethyl-isoxazol-5(4H)-one	113	C5H7N02	015731-93-8	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
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Misc :
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Instrument :
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ClientSampleId :
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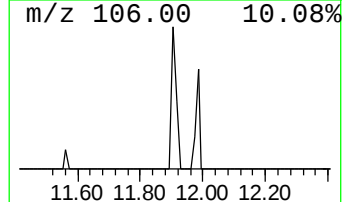
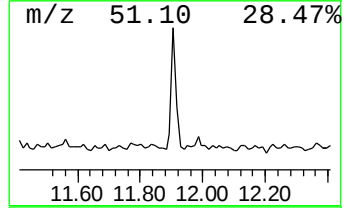
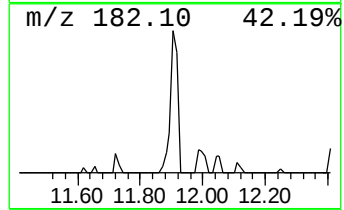
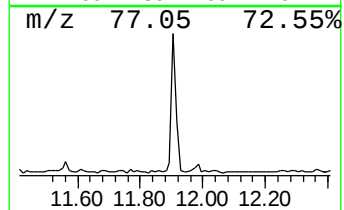
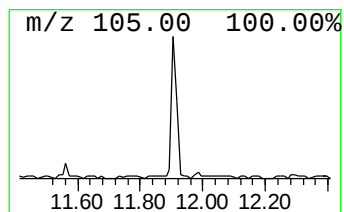
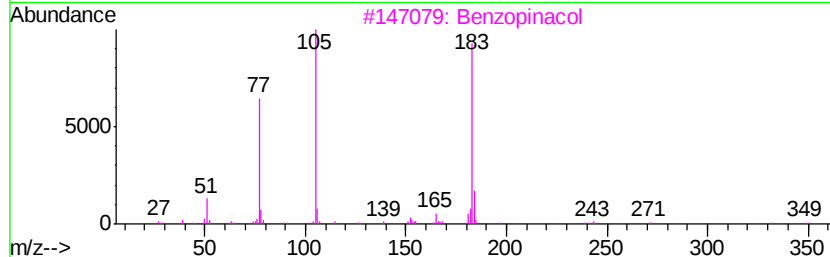
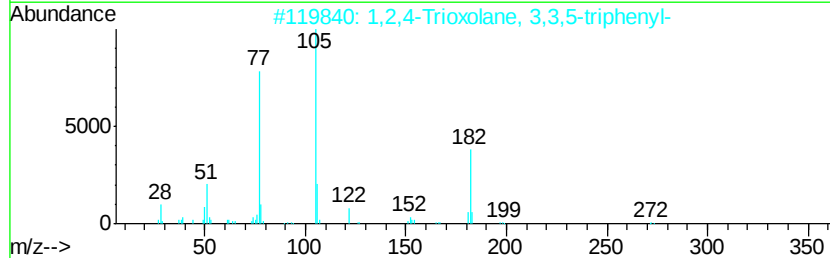
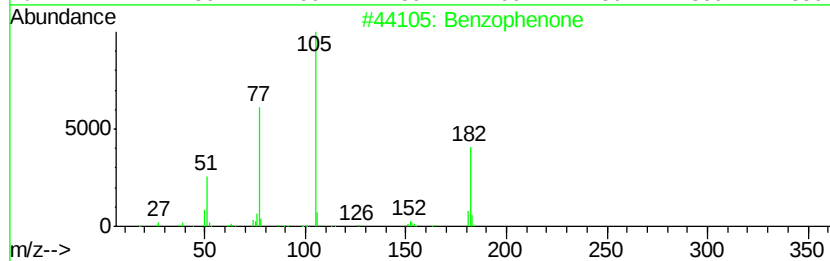
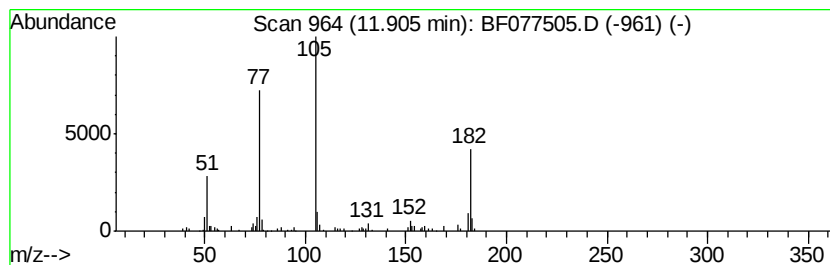
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	2.87 ng	118764	Acenaphthene-d10	11.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	64
3			Benzopinacol	366	C26H22O2	000464-72-2	47
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	43
5			Methyl benzillate	242	C15H14O3	000076-89-1	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022715\
Data File : BF077505.D
Acq On : 27 Feb 2015 14:35
Operator : TP/IZ
Sample : G1377-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.72	12.3	ng	273660	1	7.26	443886	20.0
2-Pentanone, 4-hy...	5.04	4.8	ng	105968	1	7.26	443886	20.0
unknown6.97	6.97	51.2	ng	1136200	1	7.26	443886	20.0
unknown7.47	7.47	3.1	ng	68138	1	7.26	443886	20.0
1H-Indene, 2,3-di...	8.43	2.2	ng	73632	2	8.83	675670	20.0
Naphthalene, 1-me...	9.72	2.4	ng	79999	2	8.83	675670	20.0
Naphthalene, 2-me...	9.84	8.8	ng	296574	2	8.83	675670	20.0
Naphthalene, 2,7-...	10.50	2.4	ng	101010	3	11.01	826595	20.0
Naphthalene, 2,3-...	10.59	3.8	ng	156608	3	11.01	826595	20.0
unknown11.54	11.54	2.0	ng	84099	3	11.01	826595	20.0
Benzophenone	11.91	2.9	ng	118764	3	11.01	826595	20.0