

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
 Data File : BF137332.D
 Acq On : 07 Feb 2024 21:02
 Operator : CG\JU
 Sample : P1372-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CG-MW-17-ch4

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF137332.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.875	298	304	310	rBV	879289	1174890	2.62%	0.821%
2	5.316	539	549	552	rBV	8096517	16052361	35.82%	11.222%
3	5.422	559	567	570	rBV	5149721	6057618	13.52%	4.235%
4	5.663	600	608	610	rBV	4091162	5098992	11.38%	3.565%
5	5.969	656	660	663	rBV	1179532	1074129	2.40%	0.751%
6	6.351	722	725	728	rVV	2714987	2416254	5.39%	1.689%
7	6.392	728	732	735	rVV	1034369	930014	2.08%	0.650%
8	6.422	735	737	743	rVV2	475977	533833	1.19%	0.373%
9	6.481	743	747	749	rVV	1706271	1463835	3.27%	1.023%
10	6.622	767	771	776	rVB3	3904933	5251766	11.72%	3.671%
11	6.851	806	810	813	rVV	2702903	2516504	5.62%	1.759%
12	6.981	826	832	835	rBV	5311041	5604845	12.51%	3.918%
13	7.063	840	846	849	rBV	5779533	6934184	15.47%	4.848%
14	7.322	885	890	893	rBV	997073	1026416	2.29%	0.718%
15	7.357	893	896	900	rVB	1318613	1309942	2.92%	0.916%
16	7.581	932	934	938	rVB	537275	508117	1.13%	0.355%
17	7.739	956	961	968	rVB	847254	930147	2.08%	0.650%
18	7.828	968	976	979	rBV3	3331514	5904895	13.18%	4.128%
19	7.869	979	983	988	rVB	3471798	5845253	13.04%	4.086%
20	7.945	993	996	1003	rVB	534964	570391	1.27%	0.399%
21	8.169	1013	1034	1036	rBV2	9701681	44815094	100.00%	31.329%
22	8.192	1036	1038	1040	rVB	4545904	3137681	7.00%	2.193%
23	8.733	1128	1130	1135	rVV	746162	677295	1.51%	0.473%
24	8.786	1135	1139	1143	rVV	1258198	1160853	2.59%	0.812%
25	8.898	1150	1158	1161	rVV	5991443	8630934	19.26%	6.034%
26	9.145	1196	1200	1203	rBV	2366134	1951819	4.36%	1.364%
27	9.245	1214	1217	1224	rVB	921355	716545	1.60%	0.501%
28	9.410	1238	1245	1248	rVV	604116	781561	1.74%	0.546%
29	9.480	1253	1257	1260	rVV	1023051	1040714	2.32%	0.728%
30	9.510	1260	1262	1266	rVB	517695	458017	1.02%	0.320%
31	9.598	1271	1277	1286	rBV	446621	514590	1.15%	0.360%
32	9.686	1286	1292	1295	rBV	1744525	1537793	3.43%	1.075%
33	9.822	1309	1315	1318	rBV2	522612	562796	1.26%	0.393%
34	9.863	1318	1322	1325	rVB	2313598	2199428	4.91%	1.538%
35	10.610	1442	1449	1464	rVB2	1270712	1374655	3.07%	0.961%
36	11.333	1570	1572	1579	rVB	518717	478892	1.07%	0.335%

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CG-MW-17-ch4

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.910	1836	1840	1850	rBV	2019538	1802034	4.02%	1.260%
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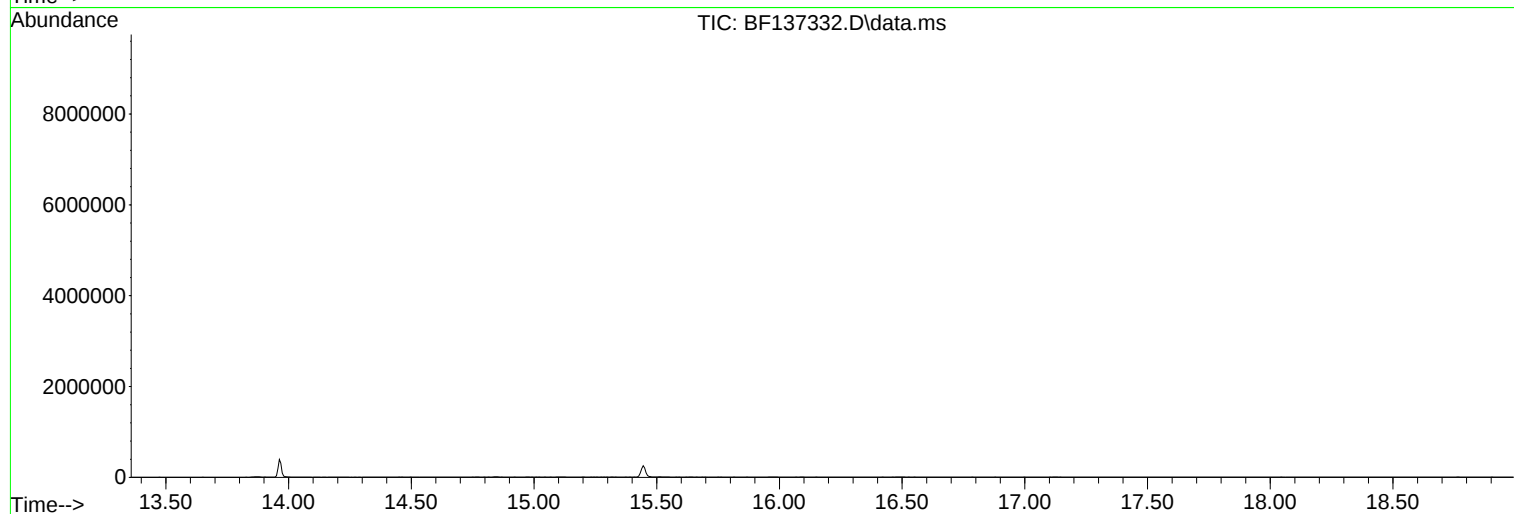
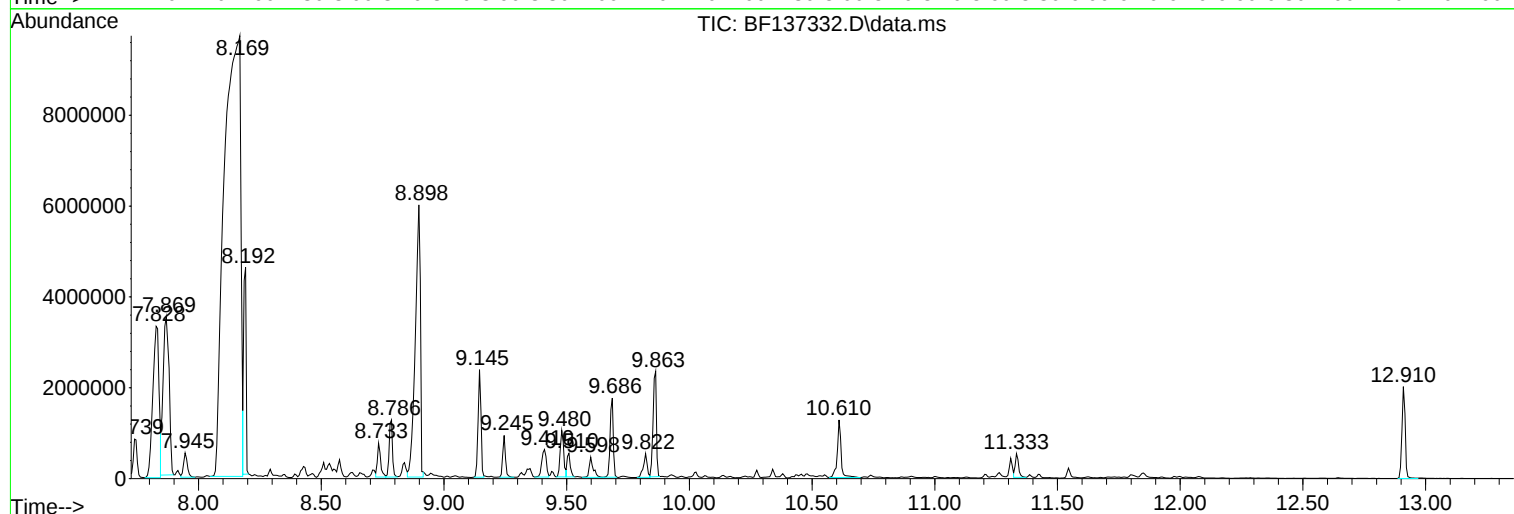
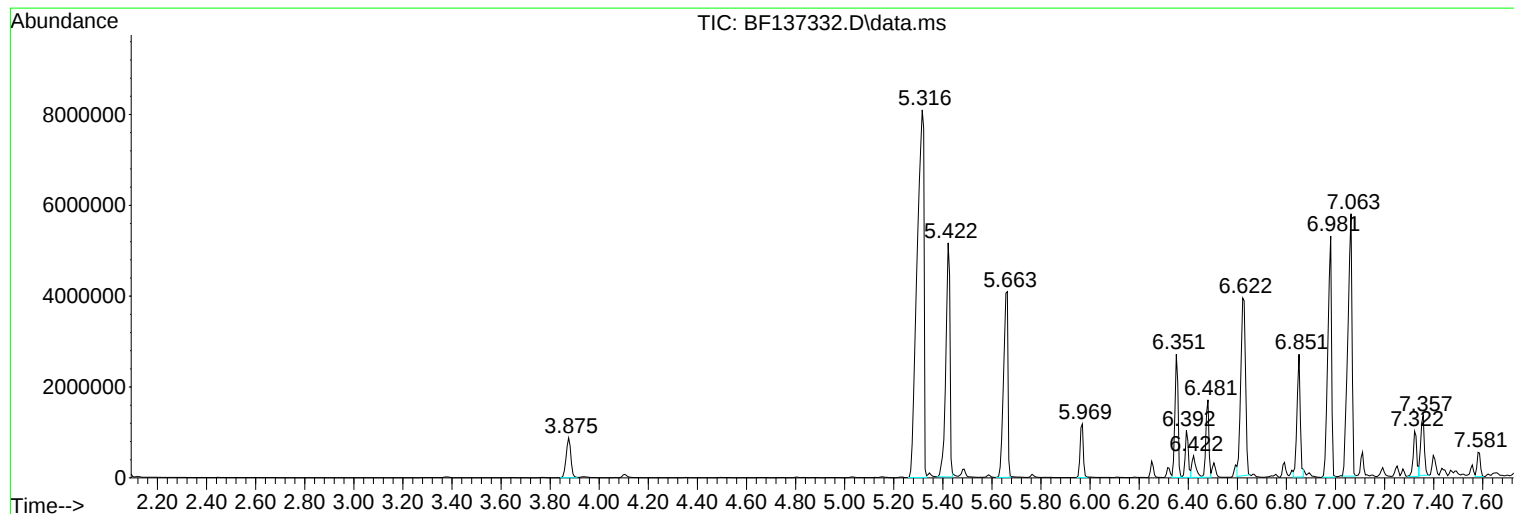
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Sample : P1372-05
Misc :
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Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

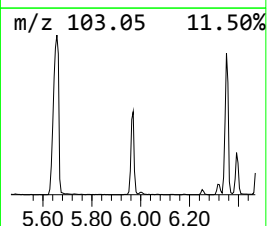
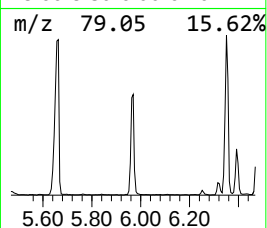
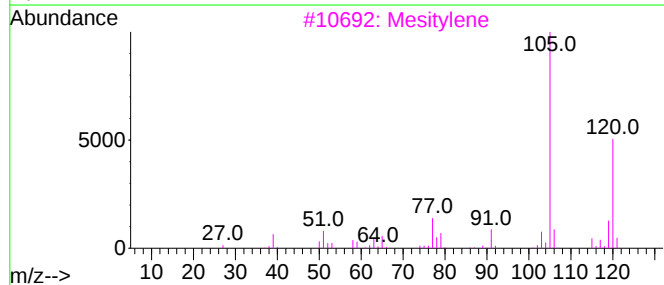
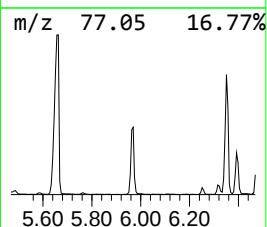
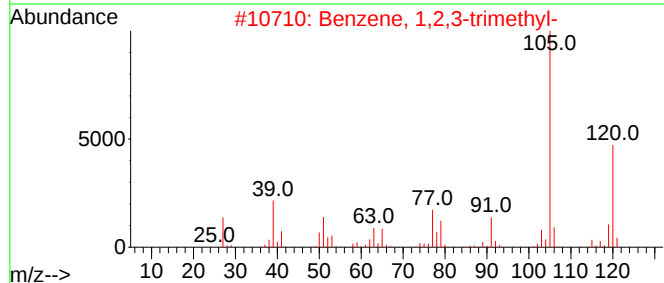
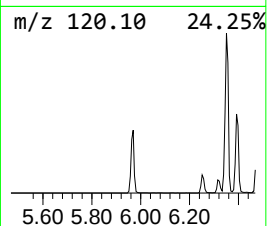
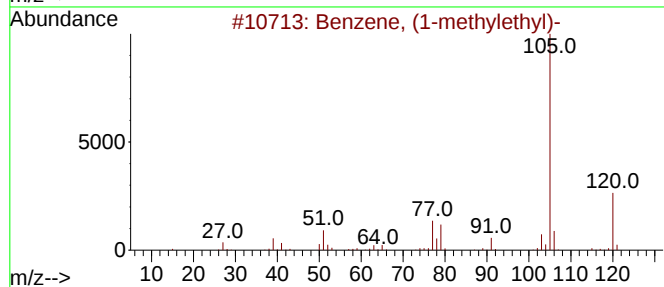
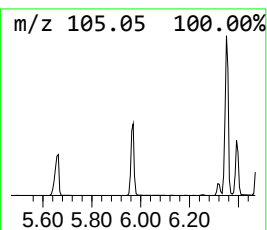
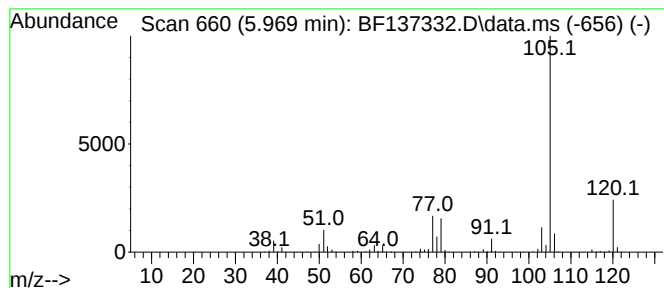
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.969	8.54 ng	1074130	1,4-Dichlorobenzene-d4	6.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	86
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	83
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

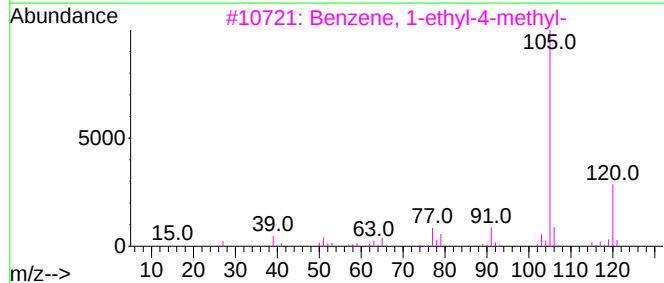
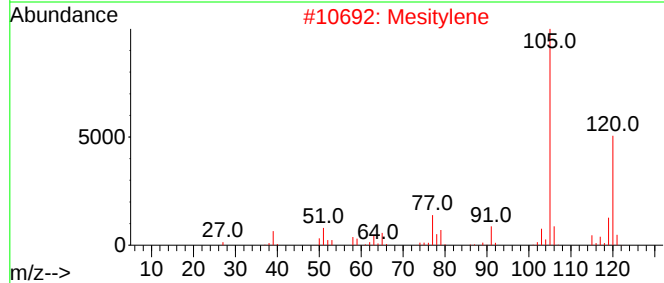
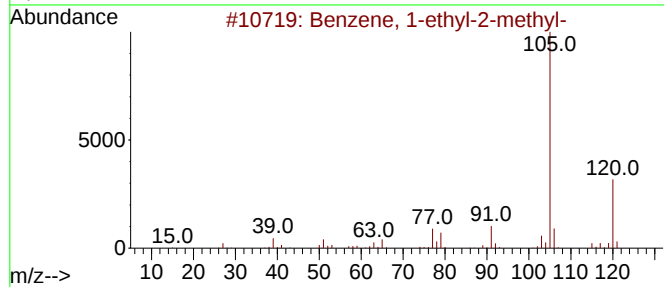
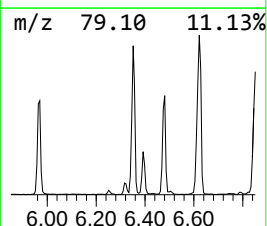
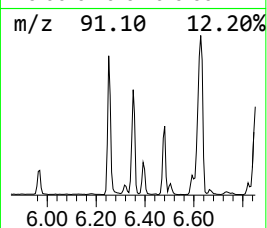
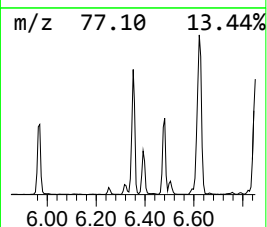
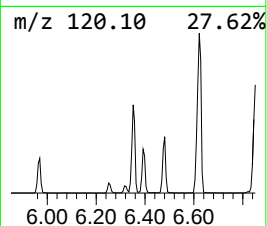
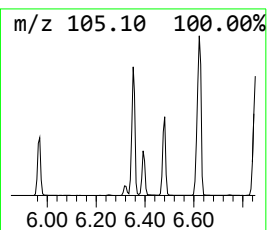
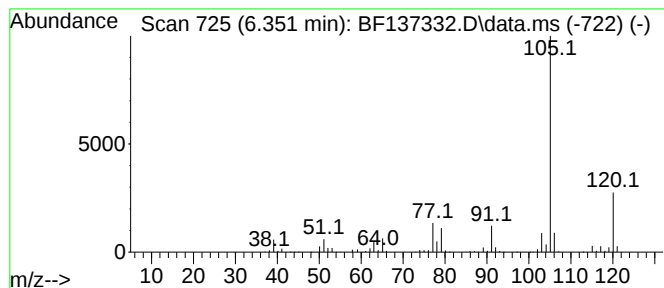
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.351	19.20 ng	2416250	1,4-Dichlorobenzene-d4	6.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

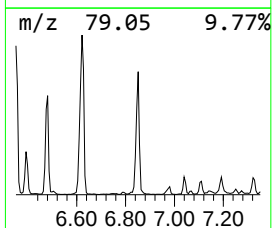
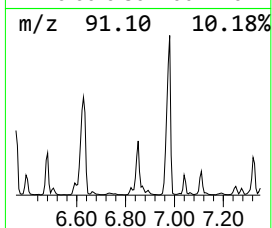
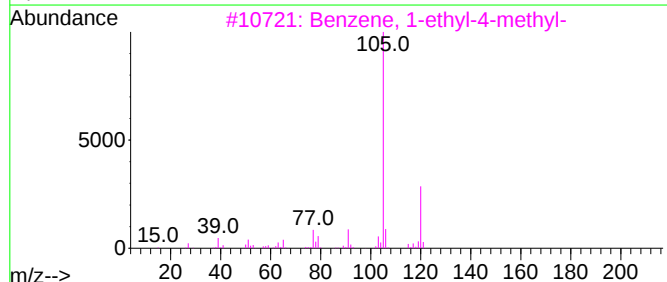
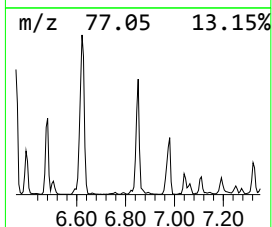
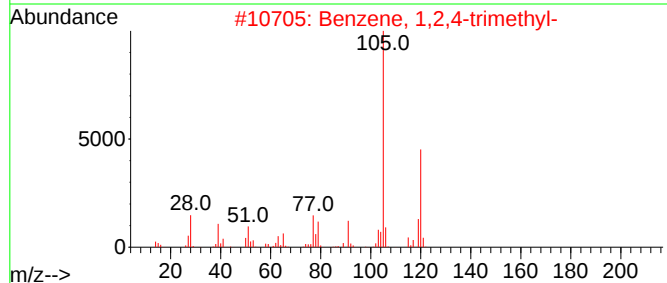
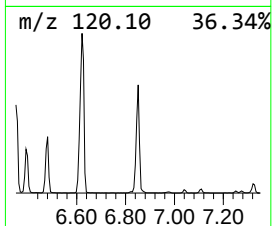
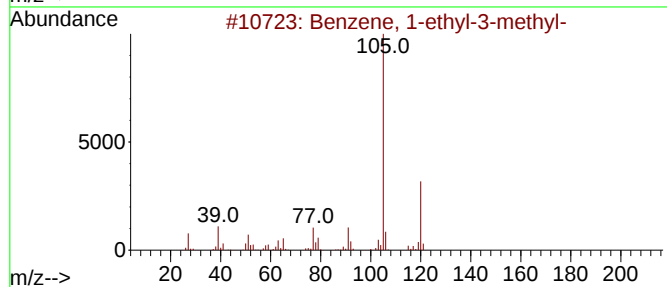
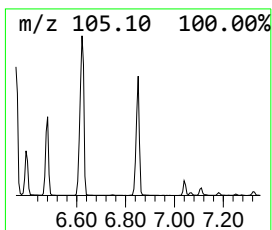
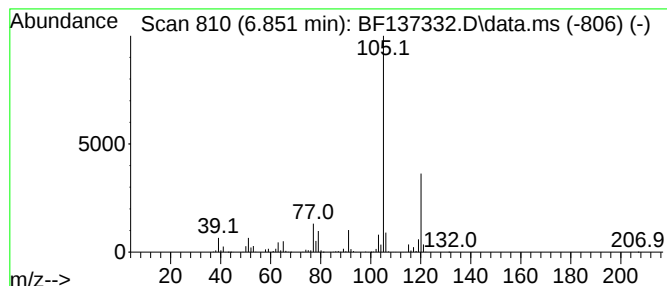
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.851	20.00 ng	2516500	1,4-Dichlorobenzene-d4	6.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Mesitylene	120	C9H12	000108-67-8	91



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

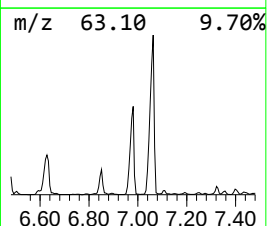
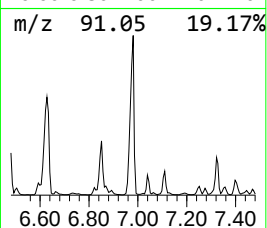
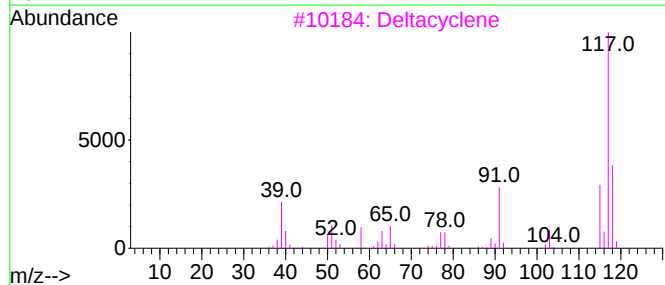
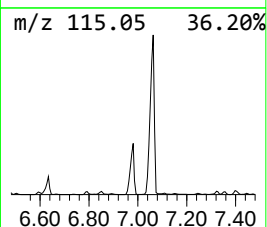
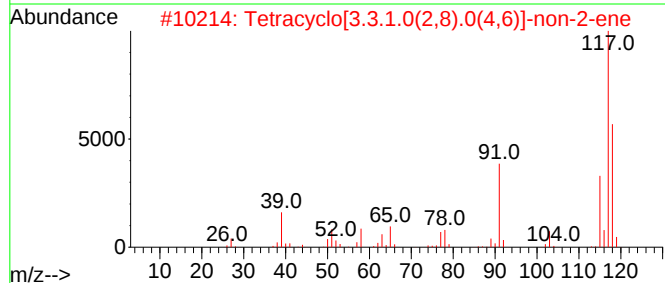
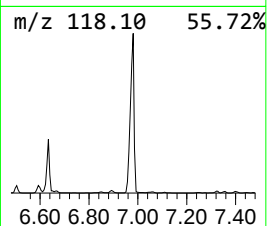
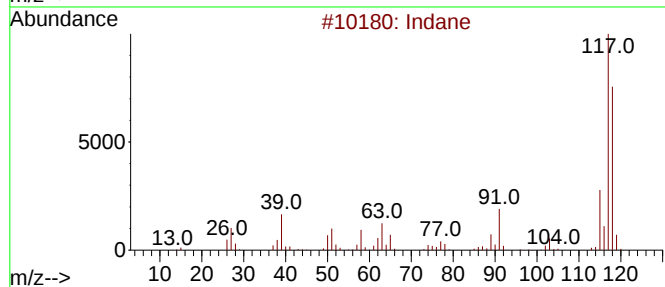
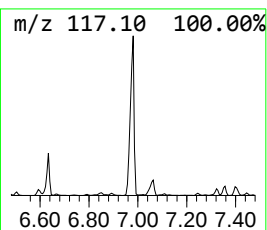
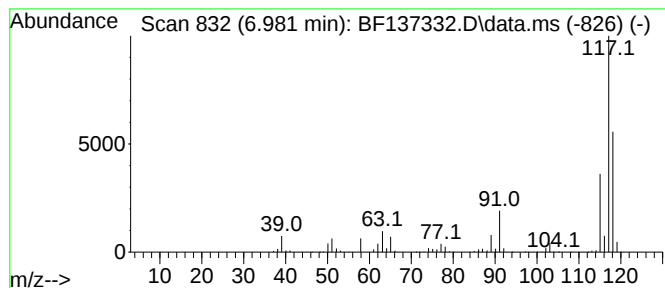
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.981	44.54 ng	5604850	1,4-Dichlorobenzene-d4	6.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	91
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	74
3			Deltacyclene	118	C9H10	007785-10-6	68
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60



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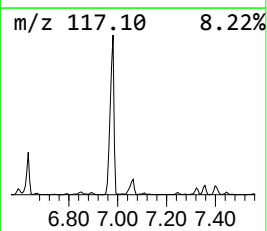
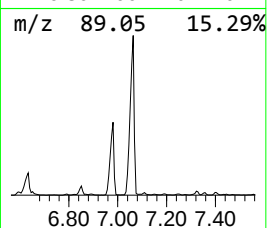
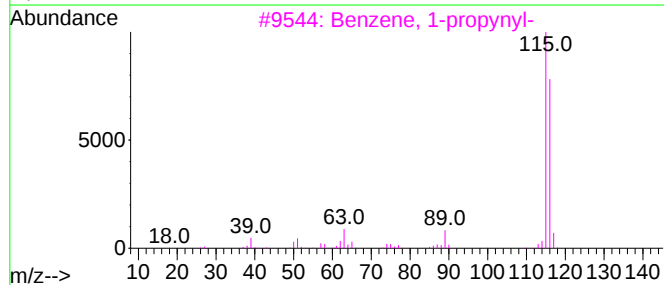
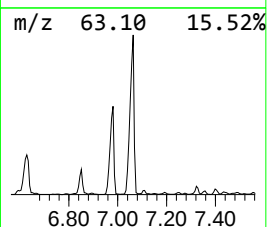
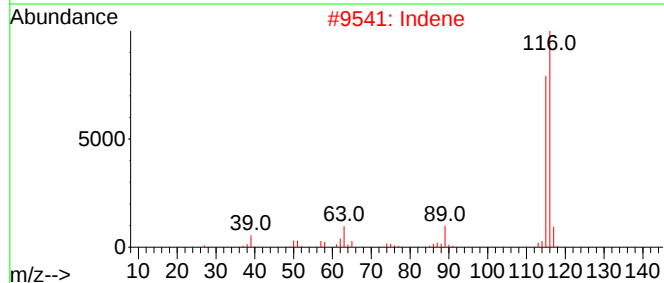
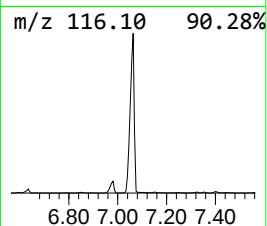
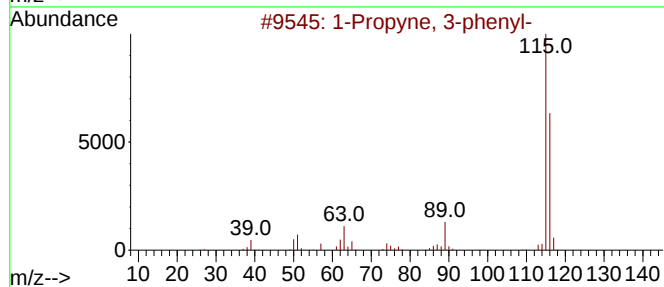
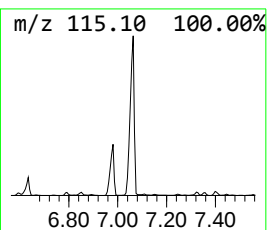
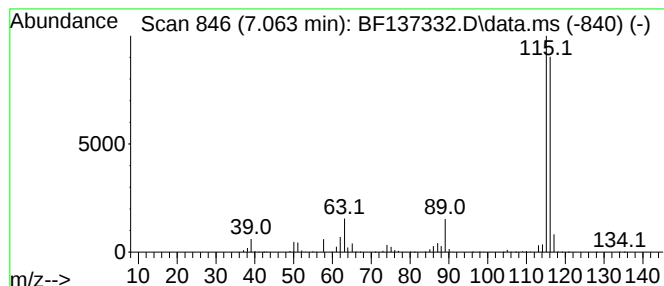
Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Propyne, 3-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
7.063	55.11 ng	6934180	1,4-Dichlorobenzene-d4		6.792	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Propyne, 3-phenyl-		116	C9H8	010147-11-2	95
2	Indene		116	C9H8	000095-13-6	94
3	Benzene, 1-propynyl-		116	C9H8	000673-32-5	91
4	3-Methylphenylacetylene		116	C9H8	000766-82-5	91
5	Benzene, 1,2-propadienyl-		116	C9H8	002327-99-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

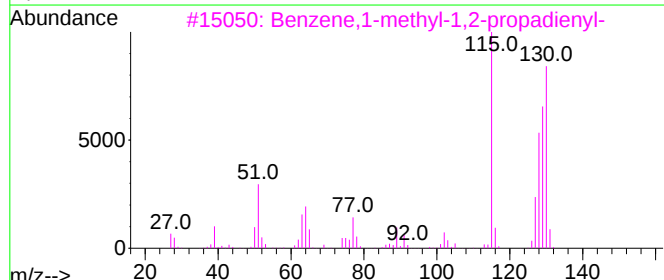
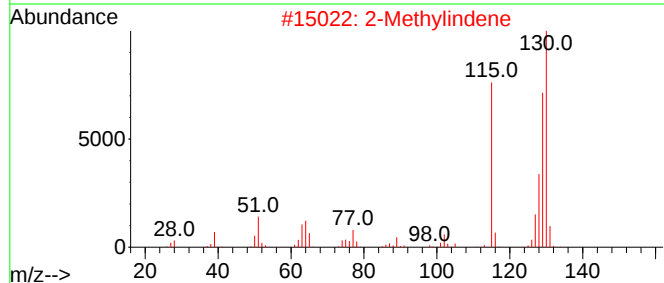
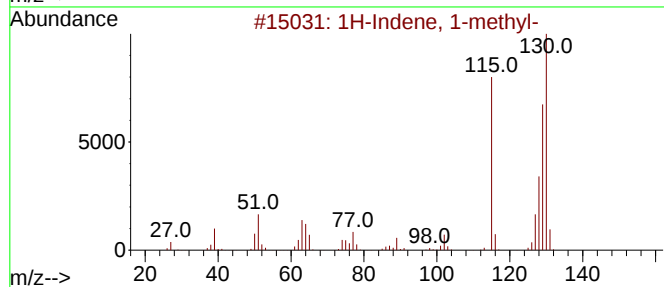
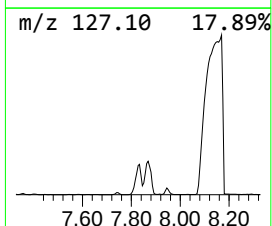
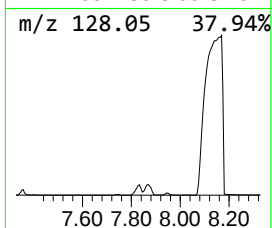
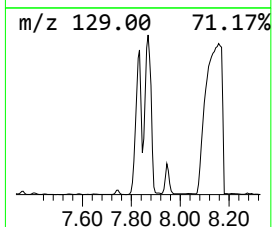
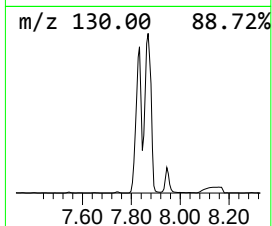
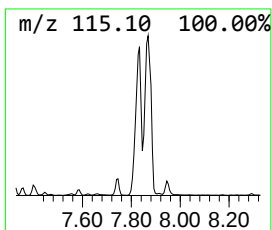
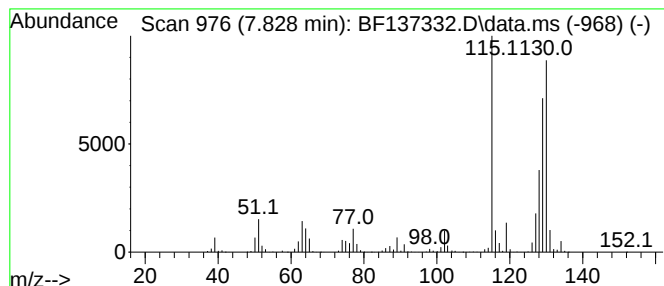
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	2.64 ng	5904900	Naphthalene-d8	8.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2		2-Methylindene	130	C10H10	002177-47-1	96
3		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	95
4		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	94
5		Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

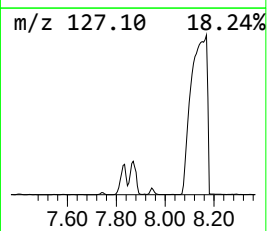
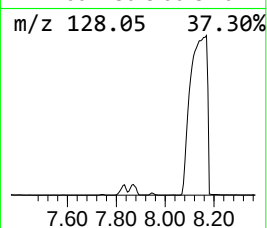
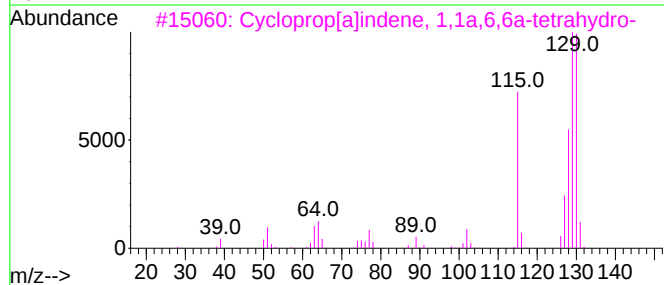
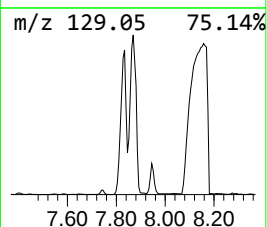
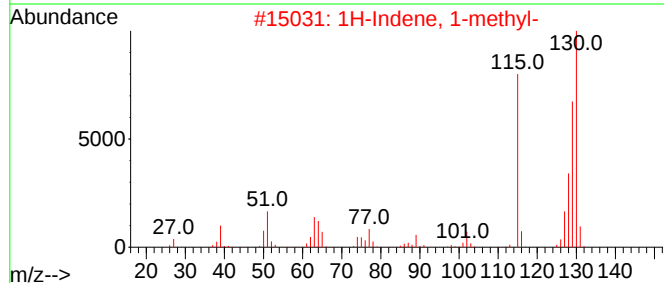
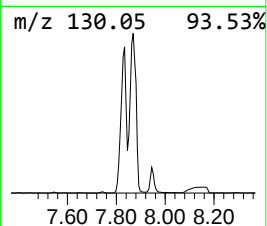
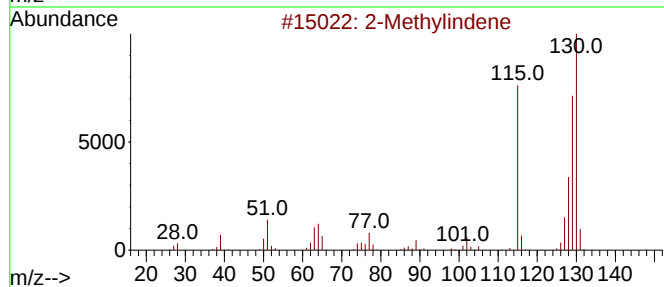
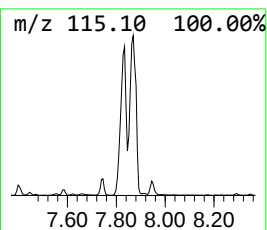
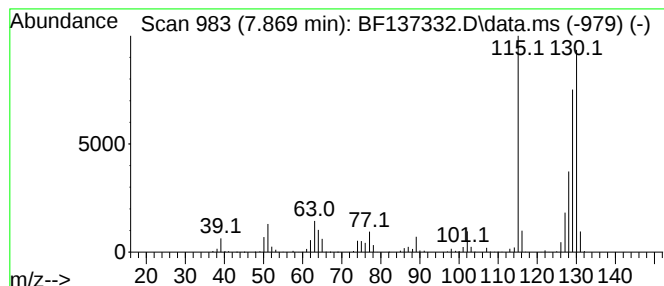
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 2-Methylindene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.869	2.61 ng	5845250	Naphthalene-d8	8.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methylindene	130	C10H10	002177-47-1	96
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
3		Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	94
4		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
5		Benzene, 1-butylnyl-	130	C10H10	000622-76-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

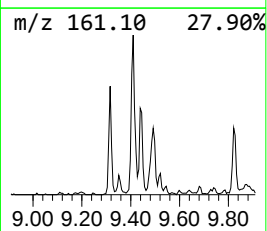
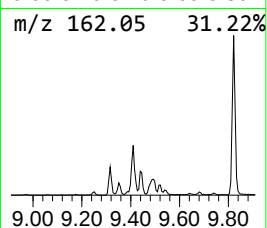
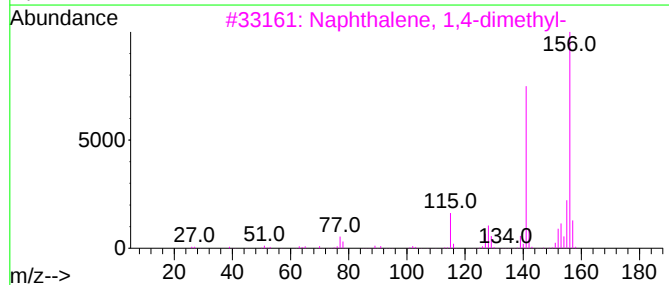
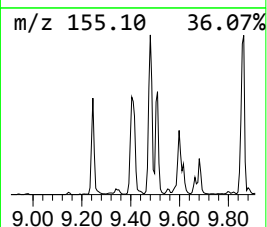
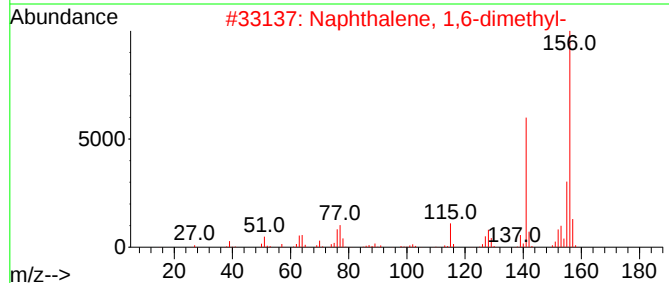
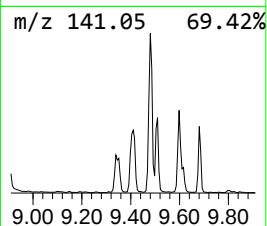
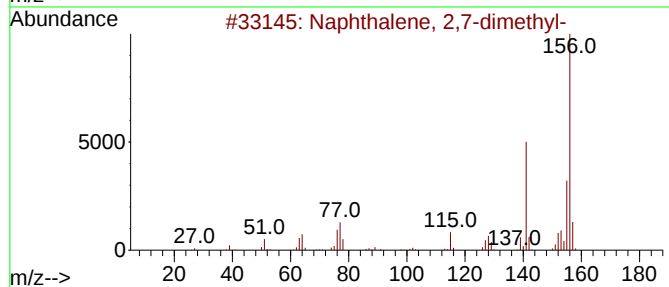
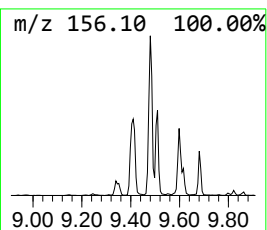
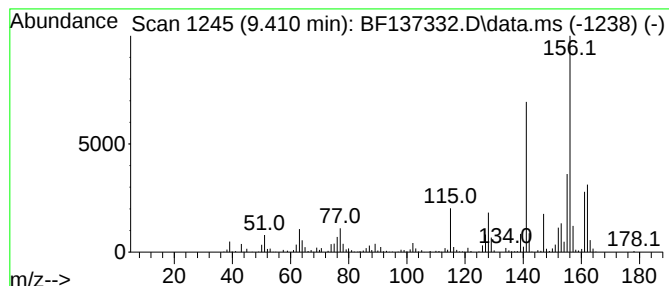
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.410	27.77 ng	781561	Acenaphthene-d10	9.822

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CG-MW-17-ch4

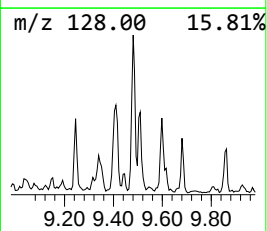
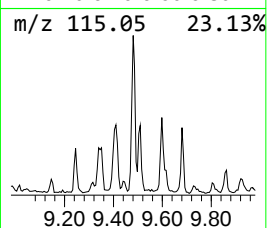
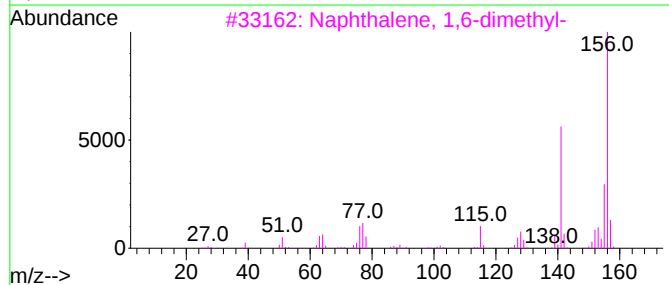
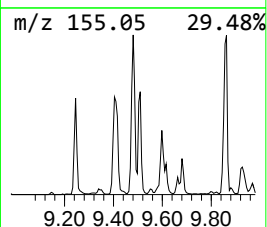
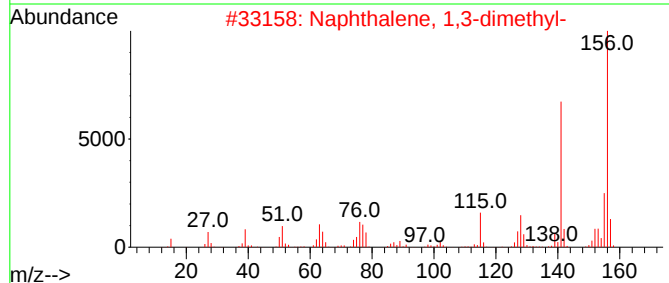
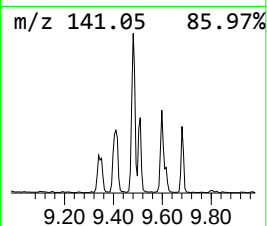
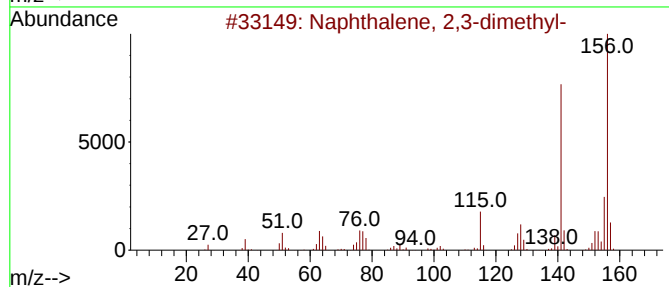
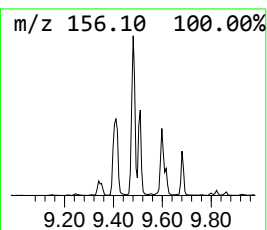
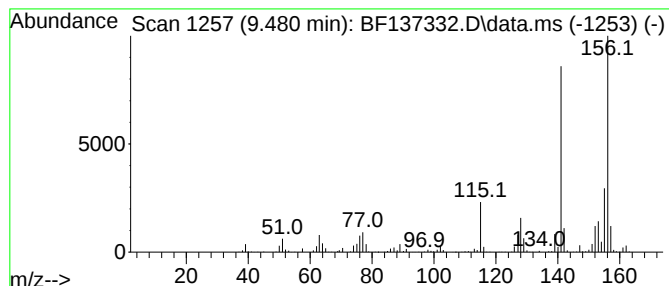
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.480	36.98 ng	1040710	Acenaphthene-d10	9.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CG-MW-17-ch4

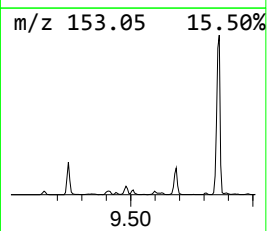
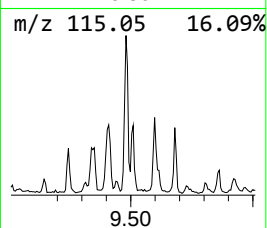
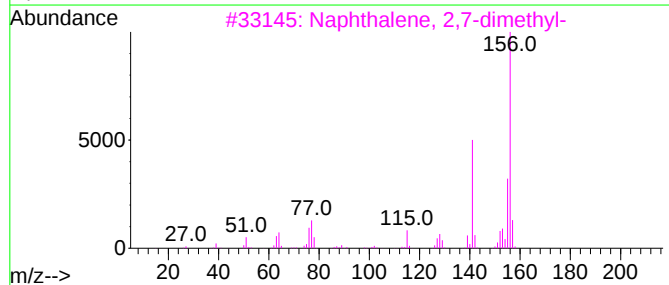
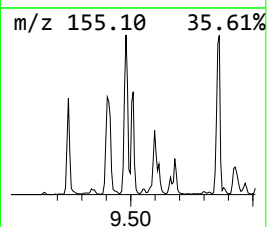
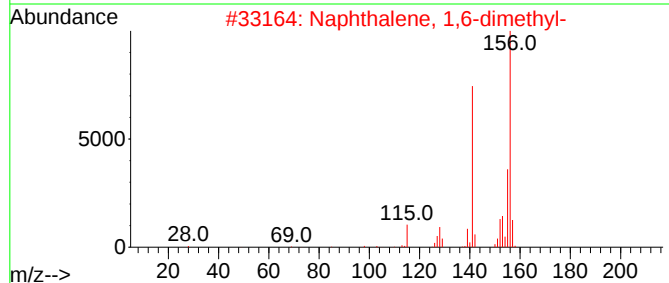
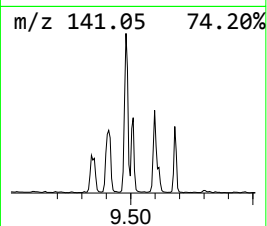
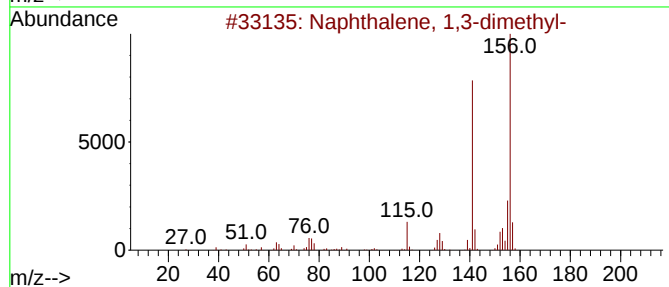
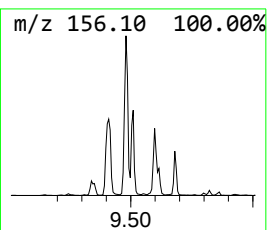
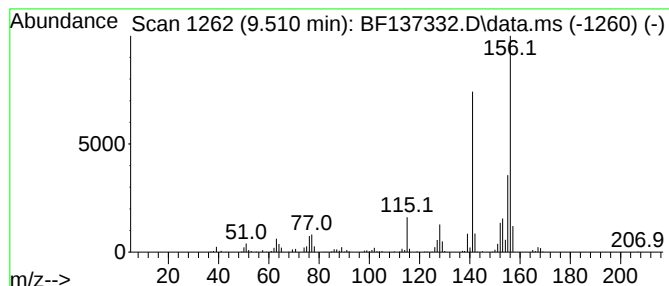
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Naphthalene, 1,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	16.28 ng	458017	Acenaphthene-d10	9.822

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CG-MW-17-ch4

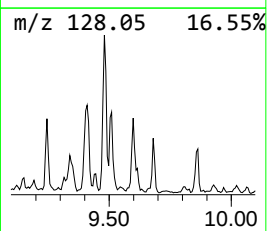
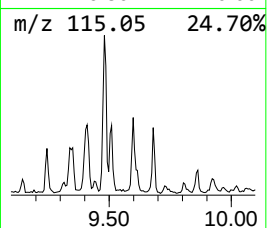
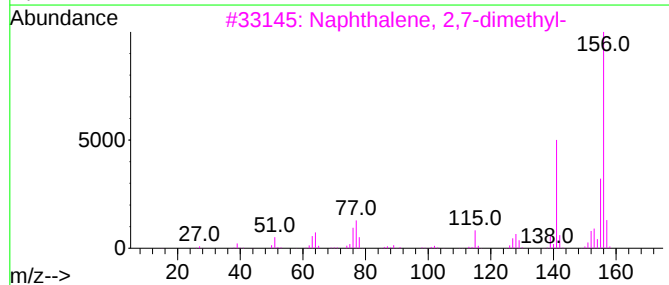
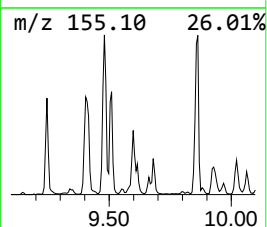
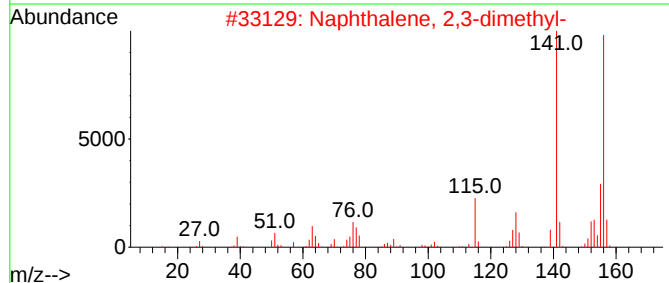
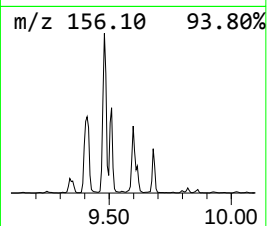
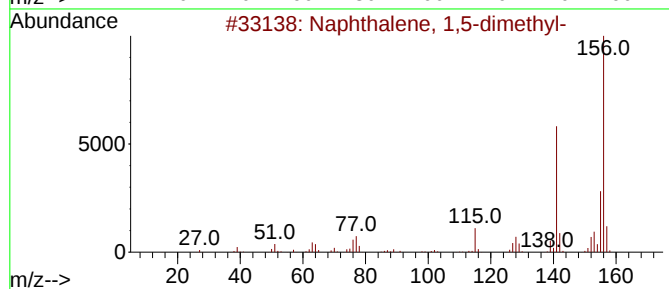
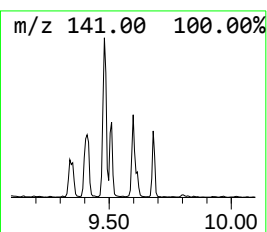
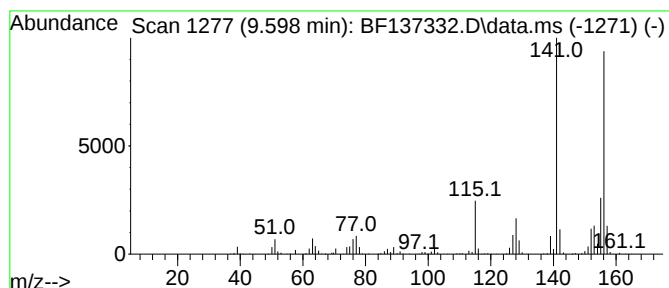
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Naphthalene, 1,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.598	18.29 ng	514590	Acenaphthene-d10	9.822

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020724\
Data File : BF137332.D
Acq On : 07 Feb 2024 21:02
Operator : CG\JU
Sample : P1372-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CG-MW-17-ch4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1-ethy...	6.351	19.2	ng	2416250	1	6.792	2516500	20.0
Benzene, 1-ethy...	6.851	20.0	ng	2516500	1	6.792	2516500	20.0
Indane	6.981	44.5	ng	5604850	1	6.792	2516500	20.0
1-Propyne, 3-ph...	7.063	55.1	ng	6934180	1	6.792	2516500	20.0
1H-Indene, 1-me...	7.828	2.6	ng	5904900	2	8.098	44815100	20.0
2-Methylindene	7.869	2.6	ng	5845250	2	8.098	44815100	20.0
Naphthalene, 2,...	9.410	27.8	ng	781561	3	9.822	562796	20.0
Naphthalene, 2,...	9.480	37.0	ng	1040710	3	9.822	562796	20.0
Naphthalene, 1,...	9.510	16.3	ng	458017	3	9.822	562796	20.0
Naphthalene, 1,...	9.598	18.3	ng	514590	3	9.822	562796	20.0