

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
 Data File : BP001373.D
 Acq On : 23 Dec 2019 18:08
 Operator : JU
 Sample : K6372-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-14

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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	6.210	696	701	710	rVB	684692	901345	1.56%	0.579%
2	7.040	833	842	847	rBV	1084057	1323504	2.30%	0.850%
3	7.481	912	917	925	rVB	1000216	1159226	2.01%	0.745%
4	7.634	938	943	949	rVB	892178	1087878	1.89%	0.699%
5	7.704	949	955	959	rBV	501469	609226	1.06%	0.391%
6	7.763	959	965	971	rVV	420061	589330	1.02%	0.379%
7	7.834	971	977	981	rVV	2271154	2688256	4.67%	1.727%
8	7.952	990	997	1003	rBV	1286546	1803731	3.13%	1.159%
9	8.069	1008	1017	1021	rVB	7118592	11373353	19.74%	7.306%
10	8.416	1064	1076	1079	rBV	2769791	3898514	6.77%	2.504%
11	8.551	1092	1099	1102	rBV	1226689	1670105	2.90%	1.073%
12	8.610	1102	1109	1112	rVB	5251032	6913395	12.00%	4.441%
13	8.722	1124	1128	1132	rVB2	776809	891011	1.55%	0.572%
14	8.834	1143	1147	1151	rBV2	924427	1125801	1.95%	0.723%
15	9.063	1178	1186	1189	rVV2	399169	639697	1.11%	0.411%
16	9.187	1197	1207	1209	rVV	1702468	2944582	5.11%	1.891%
17	9.222	1209	1213	1222	rVV2	2559528	3440417	5.97%	2.210%
18	9.551	1260	1269	1272	rBV	391658	579400	1.01%	0.372%
19	9.593	1272	1276	1280	rVB	1082467	1235578	2.15%	0.794%
20	9.840	1312	1318	1324	rVB	1676932	2140855	3.72%	1.375%
21	9.969	1328	1340	1344	rBV3	2324389	4868074	8.45%	3.127%
22	10.040	1344	1352	1357	rVV2	3090214	8332086	14.46%	5.352%
23	10.146	1366	1370	1375	rVB	915740	1151001	2.00%	0.739%
24	10.457	1400	1423	1426	rBV	15316918	57602672	100.00%	37.001%
25	11.187	1544	1547	1554	rVB	596046	841810	1.46%	0.541%
26	11.540	1597	1607	1610	rVB	5885067	9942373	17.26%	6.386%
27	11.698	1622	1634	1637	rBV	6263804	10157518	17.63%	6.525%
28	12.128	1700	1707	1711	rBV	2130784	2577859	4.48%	1.656%
29	12.281	1724	1733	1736	rBV	768779	956538	1.66%	0.614%
30	12.428	1753	1758	1768	rVB	604463	1051845	1.83%	0.676%
31	12.663	1790	1798	1802	rBV	898887	1293594	2.25%	0.831%
32	13.204	1885	1890	1895	rBV	561950	722824	1.25%	0.464%
33	13.263	1895	1900	1904	rVV	1749728	2151555	3.74%	1.382%
34	13.645	1955	1965	1975	rBV	608552	908677	1.58%	0.584%

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Instrument :
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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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.104	2039	2043	2050	rVB	548523	649682	1.13%	0.417%
36	14.504	2106	2111	2116	rVB	1373223	1756060	3.05%	1.128%
37	15.639	2301	2304	2309	rVV	818700	939563	1.63%	0.604%
38	17.892	2681	2687	2691	rBV	2468885	2759185	4.79%	1.772%

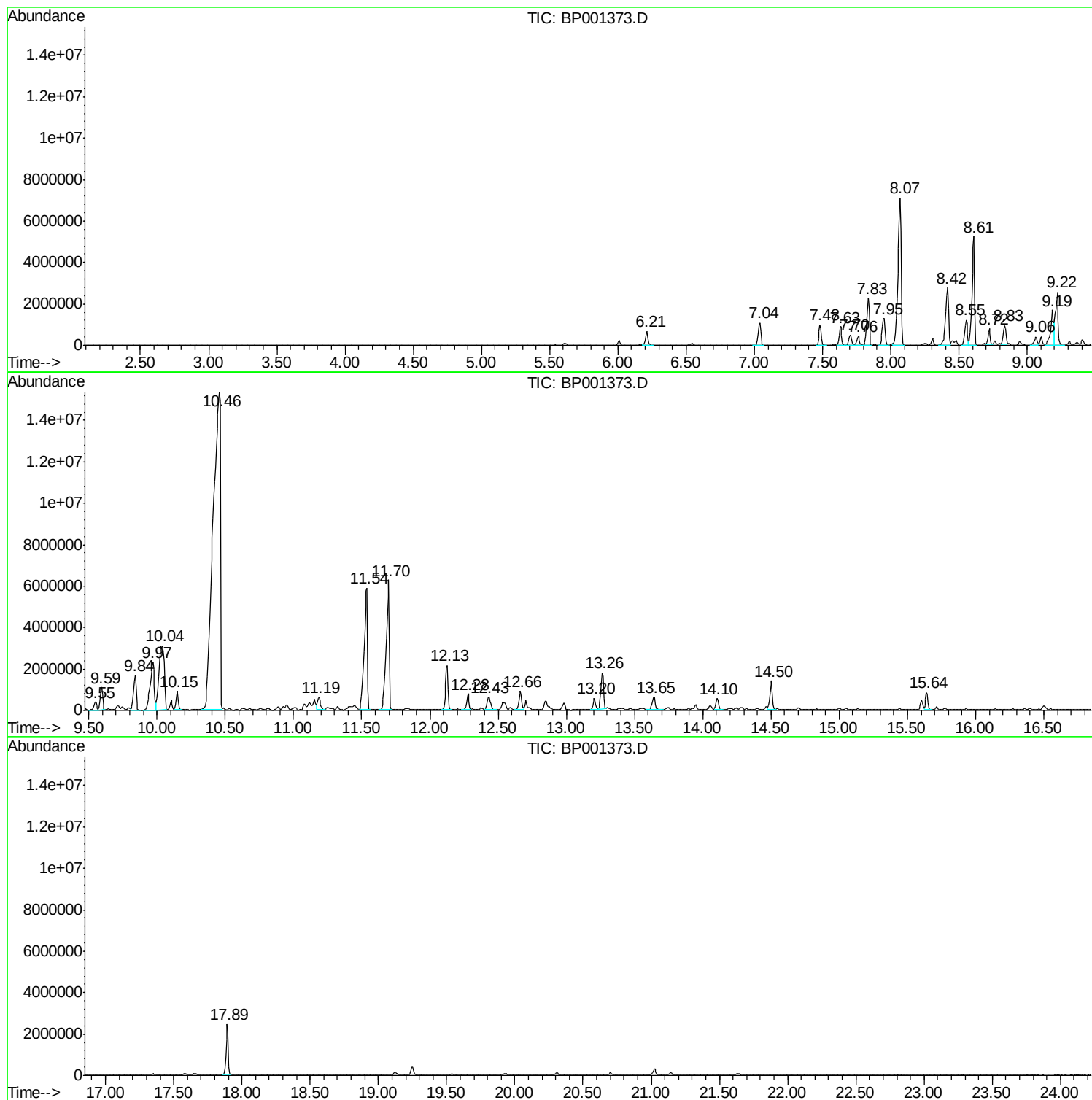
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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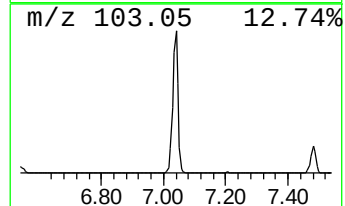
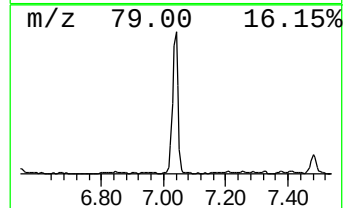
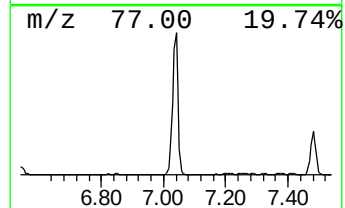
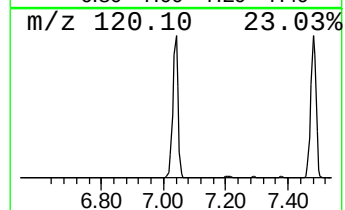
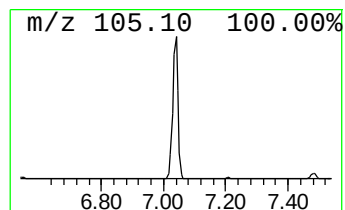
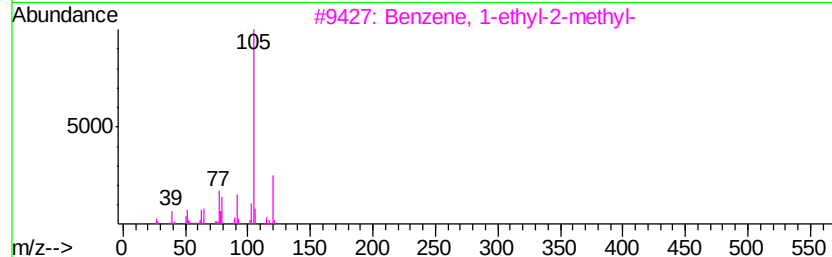
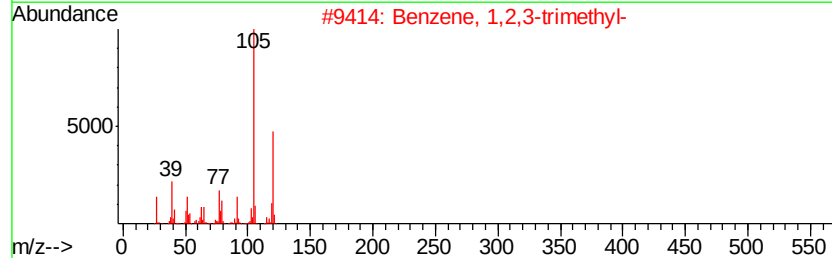
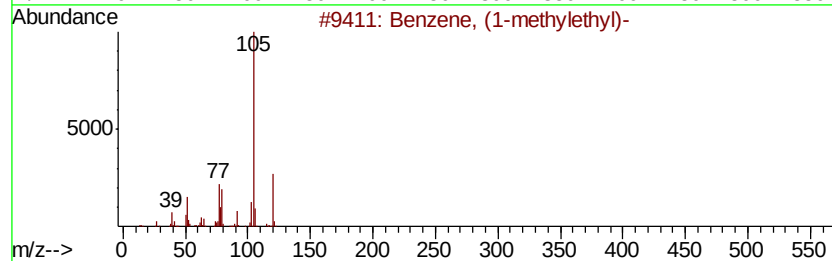
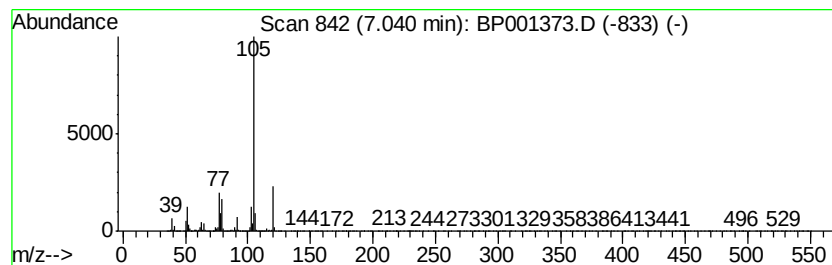
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzene, (1-methylethyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	6.79 ng	1323500	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	96
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	80
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72



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Instrument :
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ClientSampleId :
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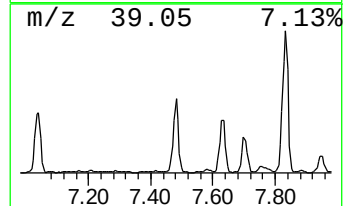
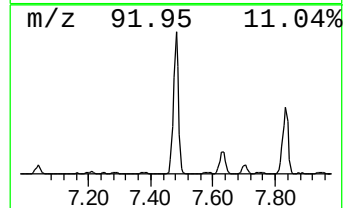
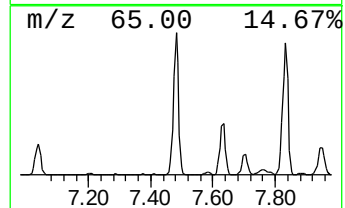
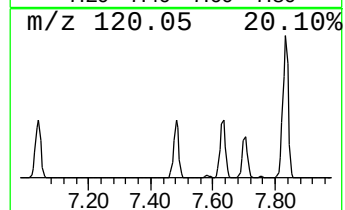
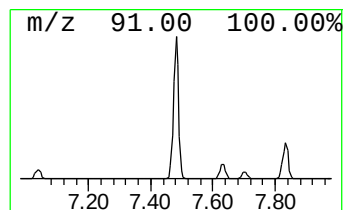
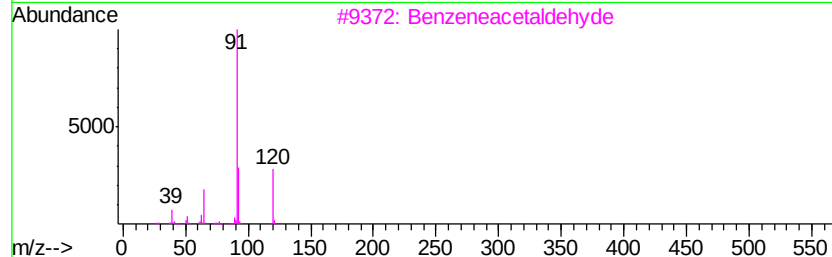
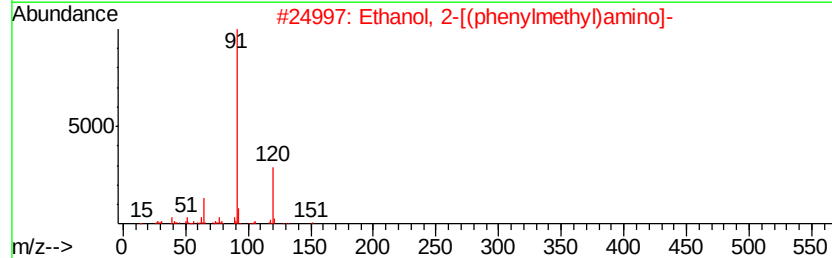
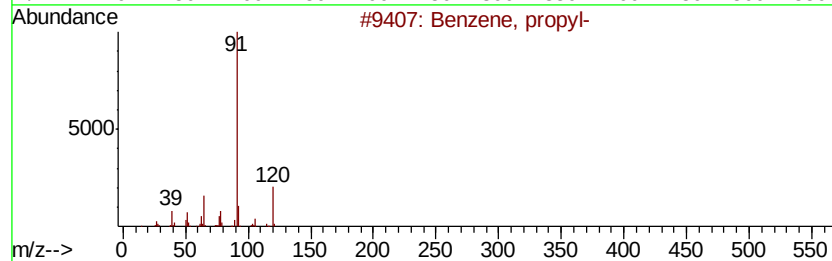
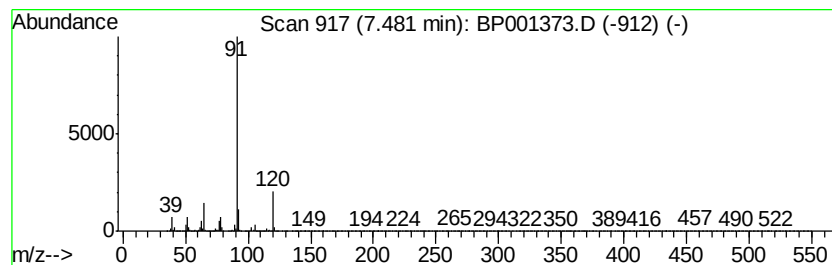
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	5.95 ng	1159230	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	95
2		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
3		Benzeneacetaldehyde	120	C8H8O	000122-78-1	62
4		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	53
5		6-Difluoromethyl-5-phenylmethane...	316	C12H10F2N2O4S	1000317-93-5	47



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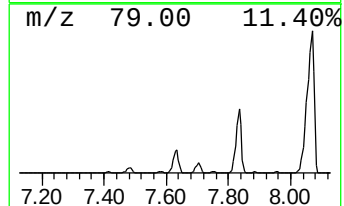
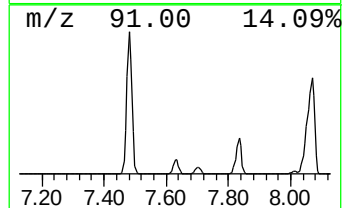
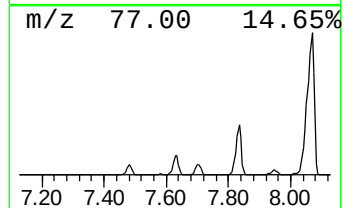
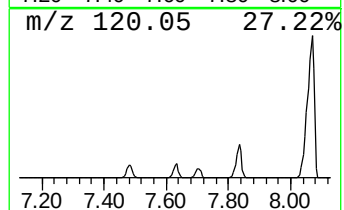
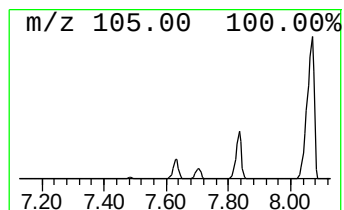
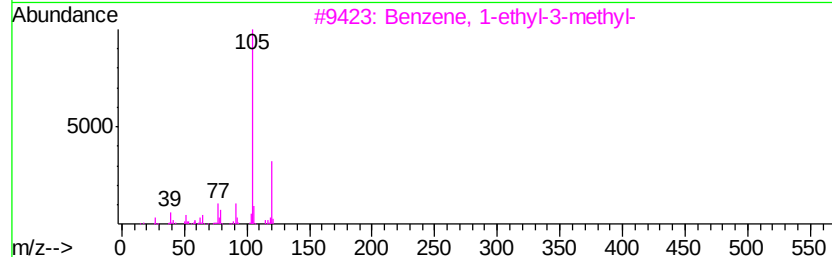
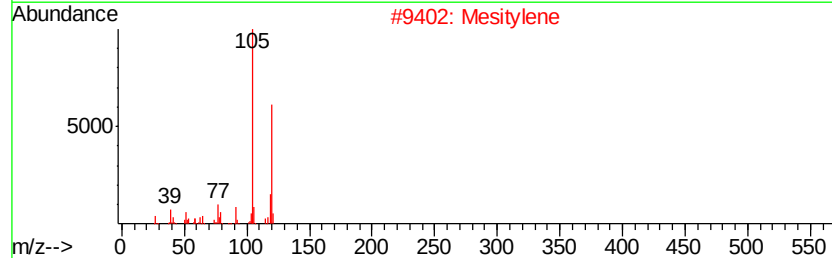
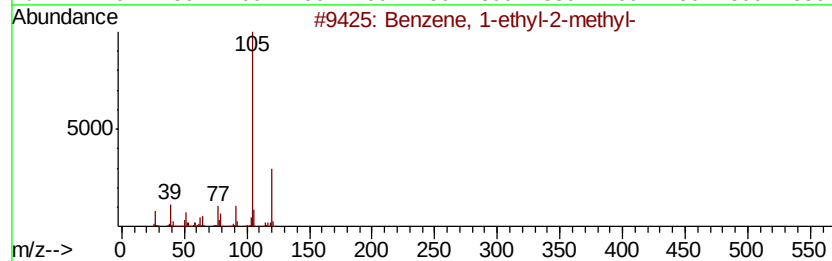
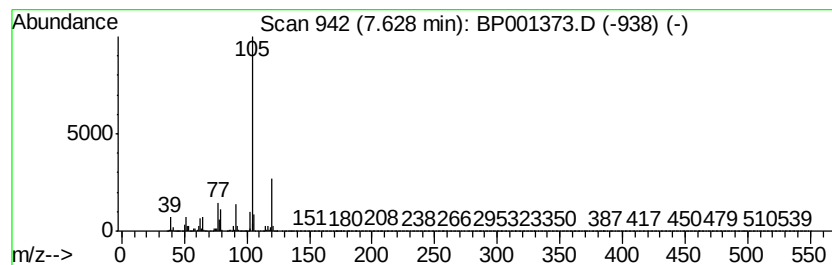
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	5.58 ng	1087880	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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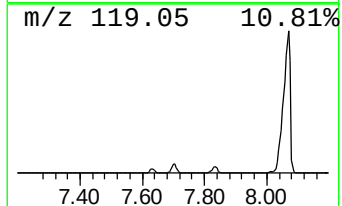
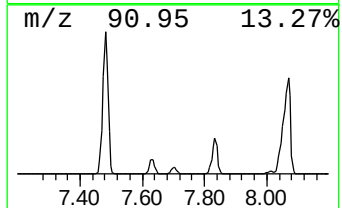
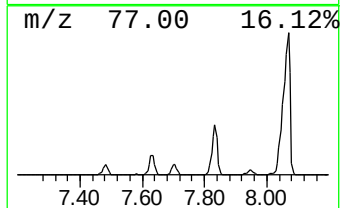
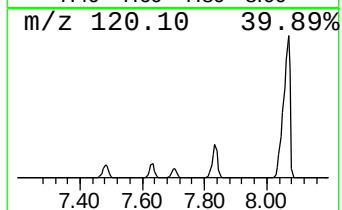
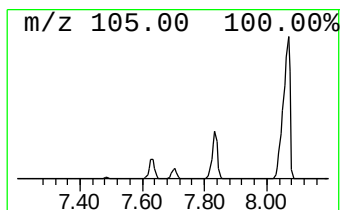
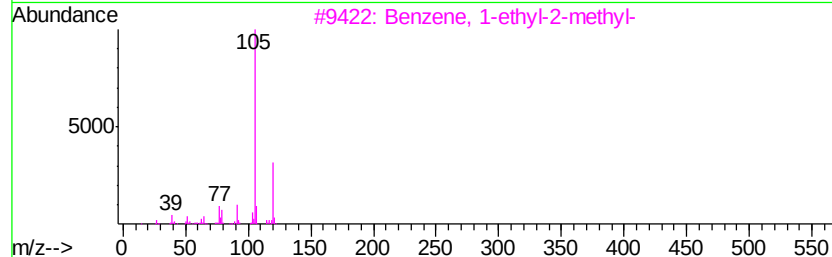
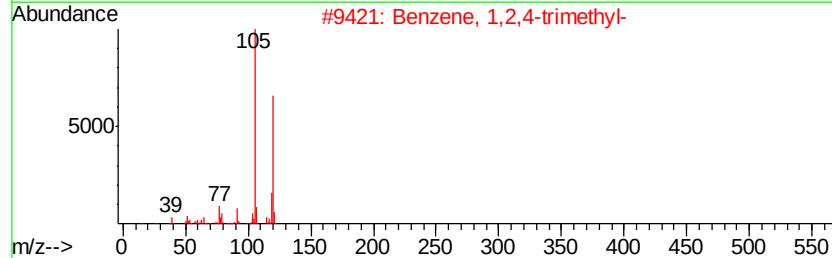
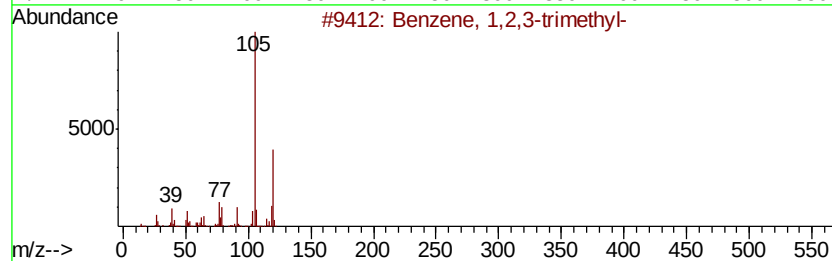
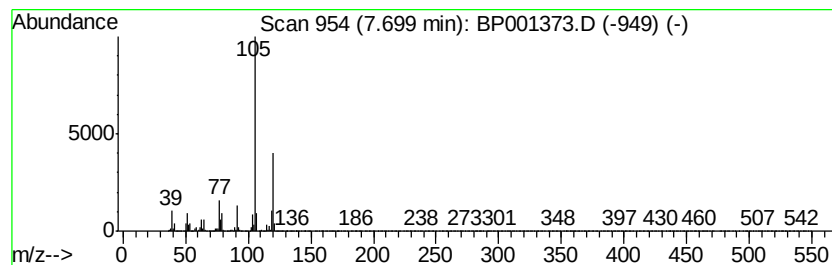
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	3.13 ng	609226	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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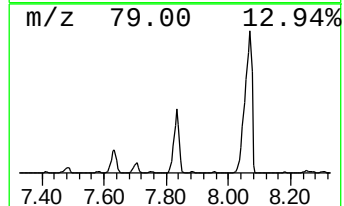
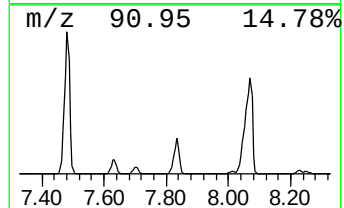
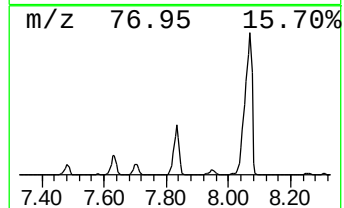
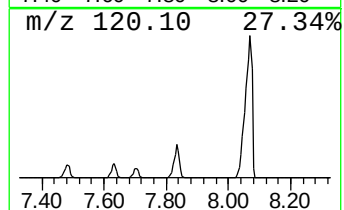
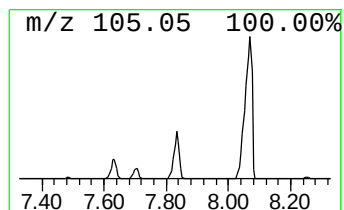
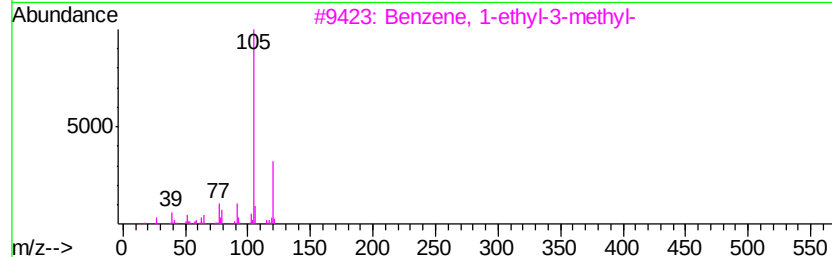
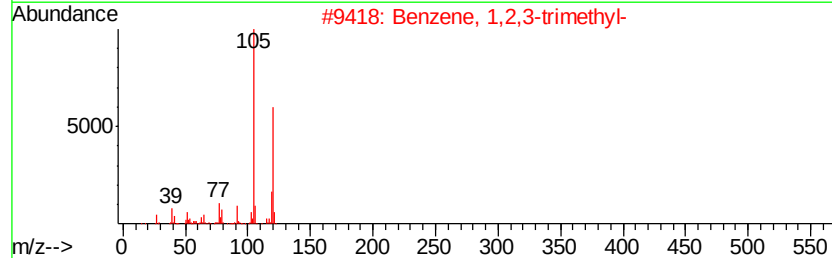
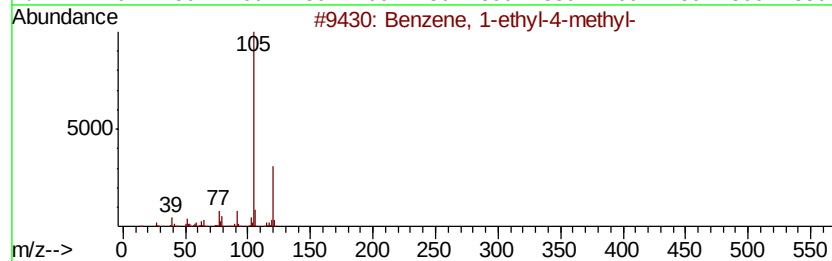
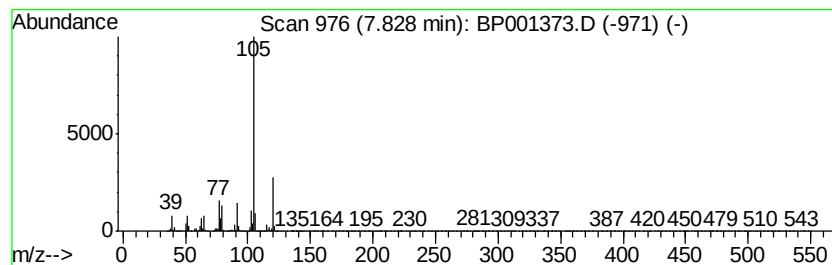
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	13.79 ng	2688260	1,4-Dichlorobenzene-d4	8.30

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-14

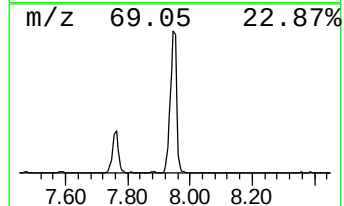
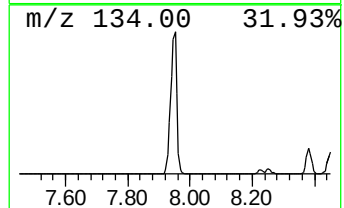
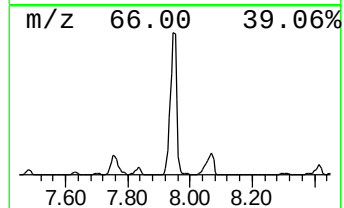
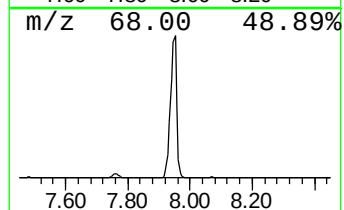
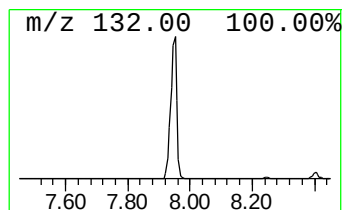
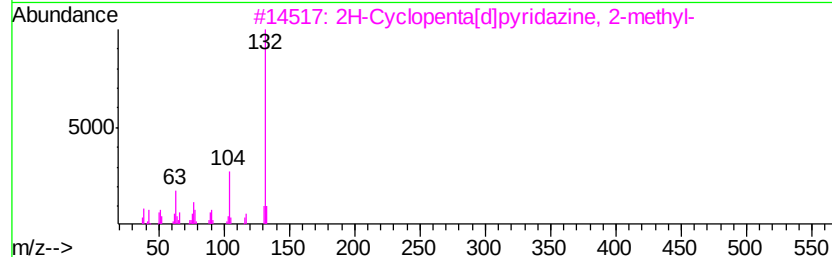
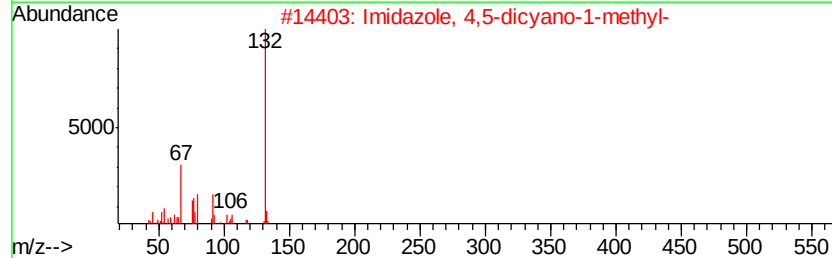
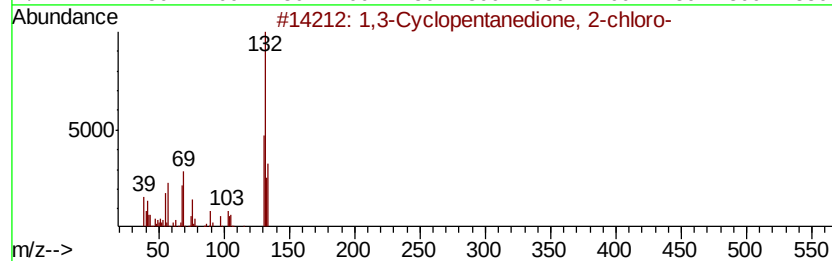
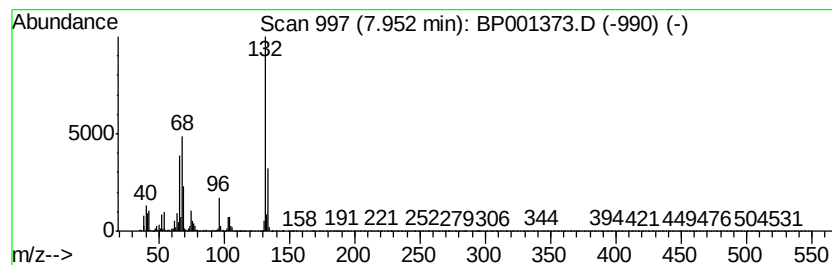
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown 7.95 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	9.25 ng	1803730	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	16
2		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	11
3		2H-Cyclopenta[d]pyridazine, 2-methyl-	132	C8H8N2	022291-85-6	10
4		Tranvlcyproamine-propionyl	189	C12H15NO	1000123-86-3	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

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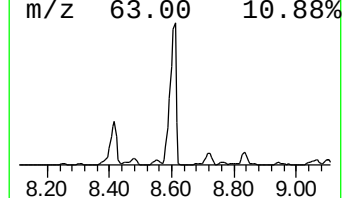
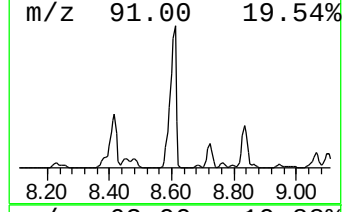
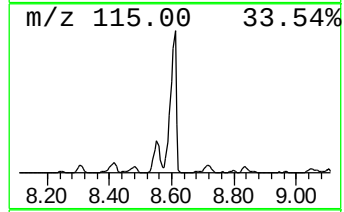
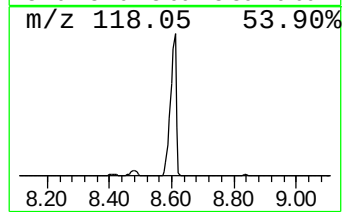
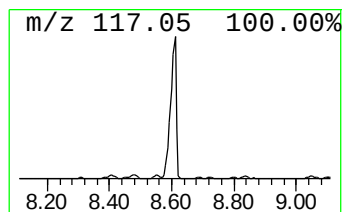
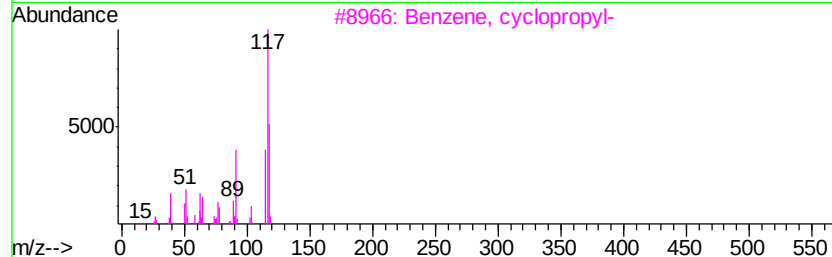
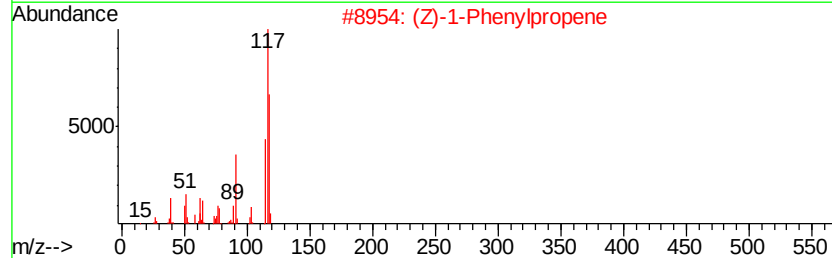
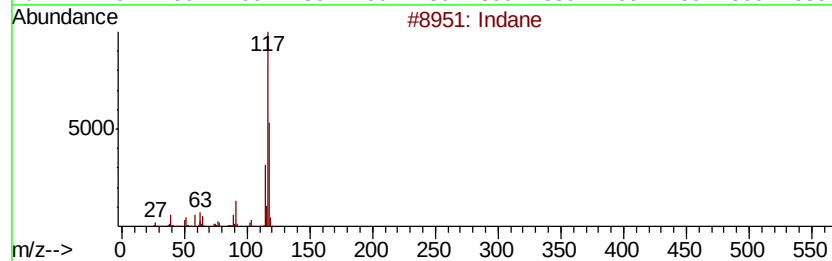
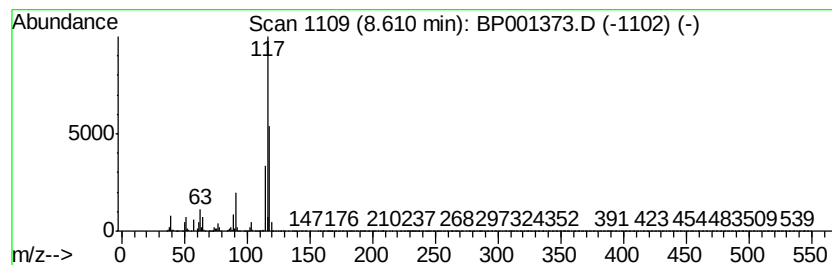
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	35.47 ng	6913400	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	87
2		(Z)-1-Phenylpropene	118	C9H10	000766-90-5	81
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4		Deltacyclene	118	C9H10	007785-10-6	76
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
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Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-14

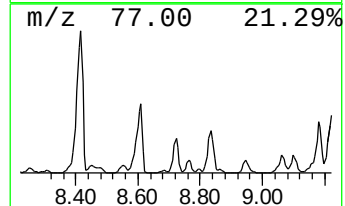
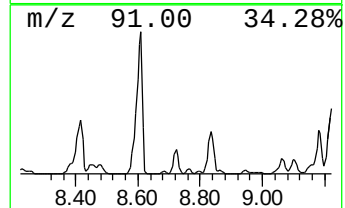
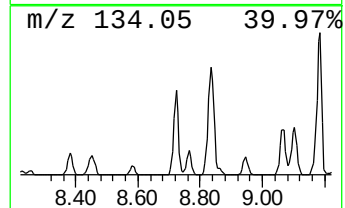
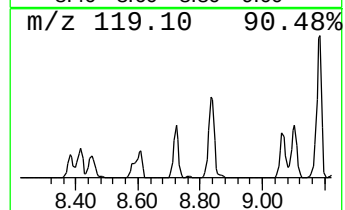
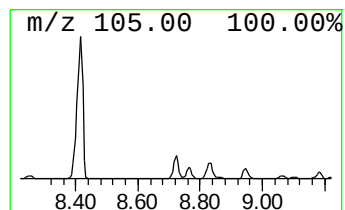
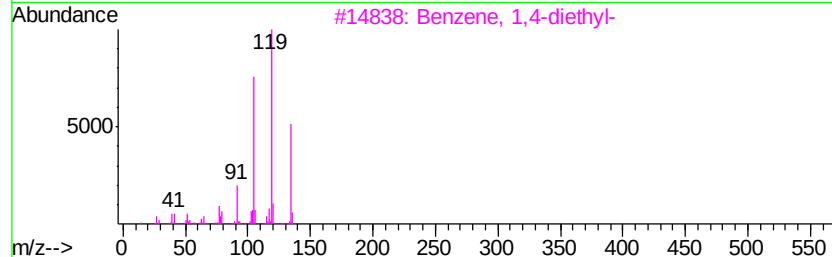
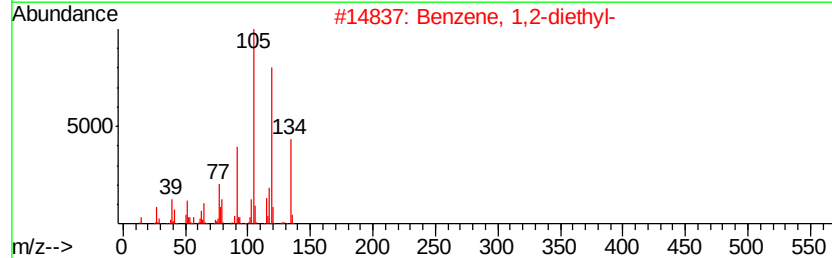
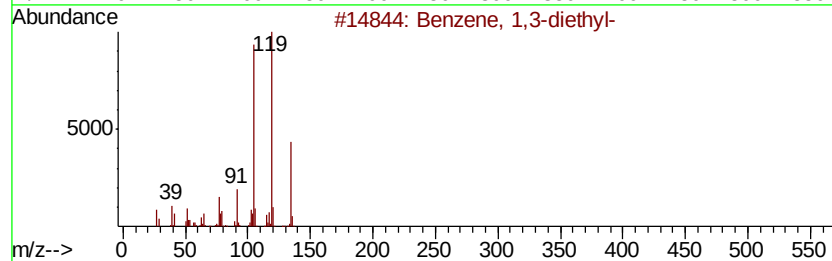
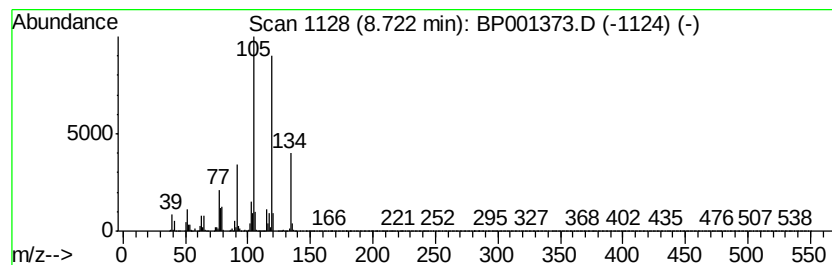
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,3-diethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	4.57 ng	891011	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

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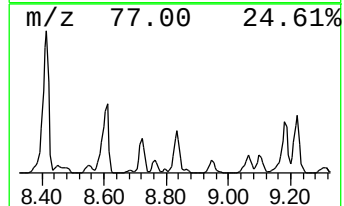
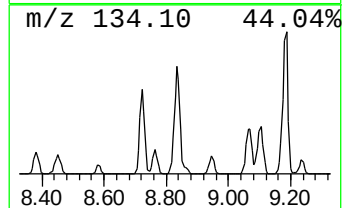
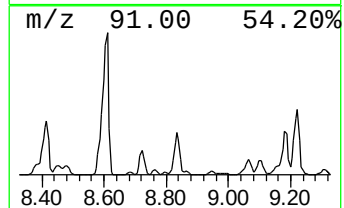
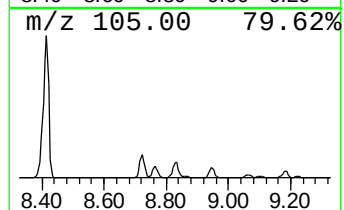
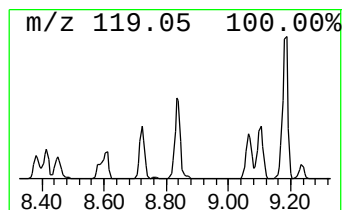
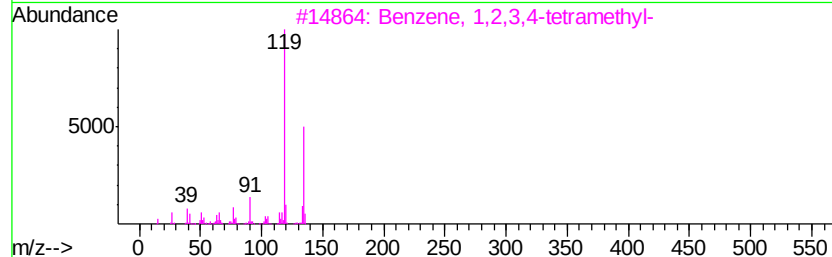
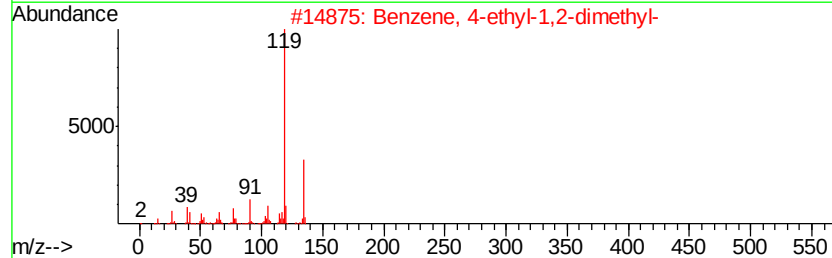
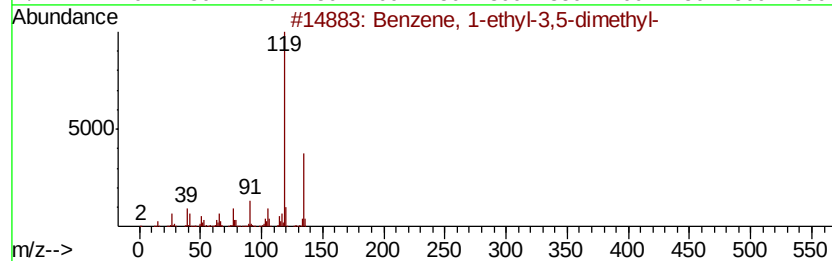
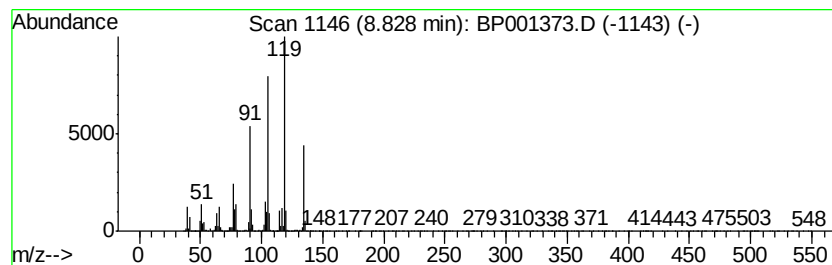
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.83	5.78 ng	1125800	1,4-Dichlorobenzene-d4	8.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
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Misc :
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Instrument :
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ClientSampleId :
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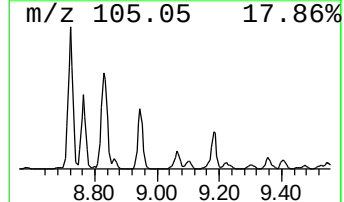
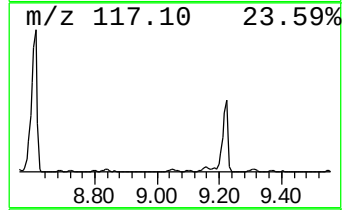
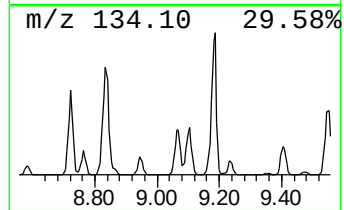
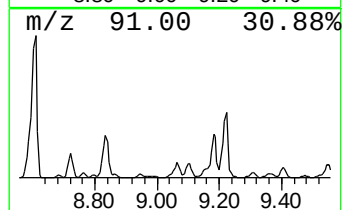
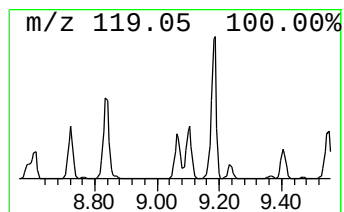
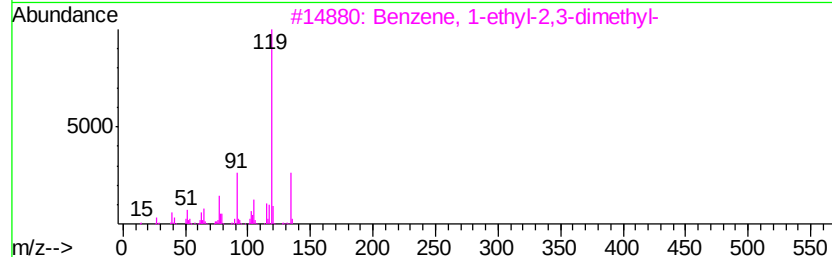
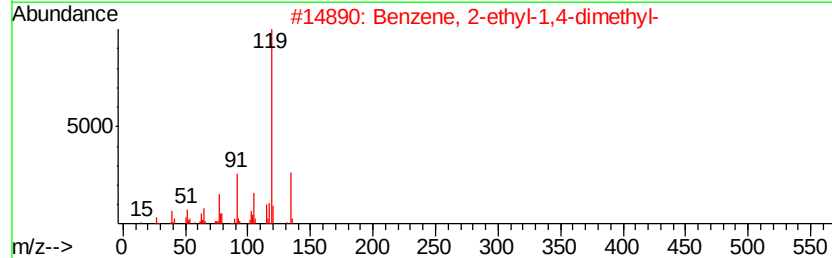
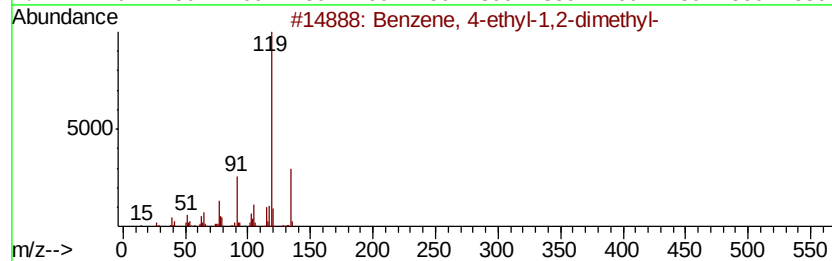
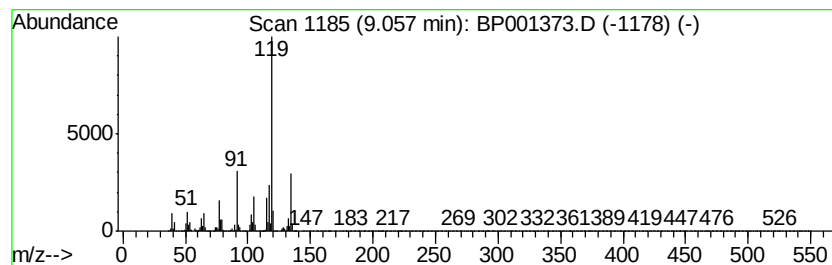
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.06	3.28 ng	639697	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
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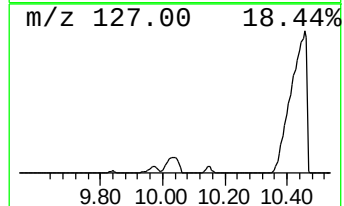
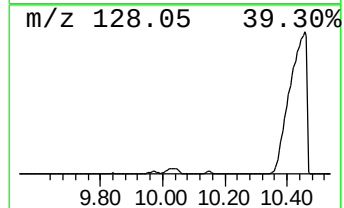
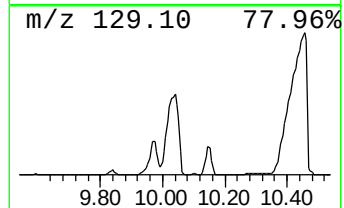
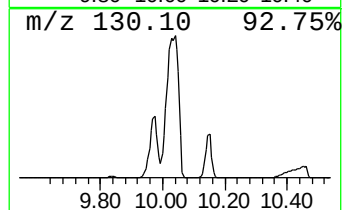
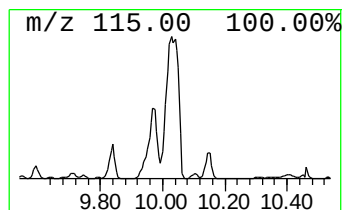
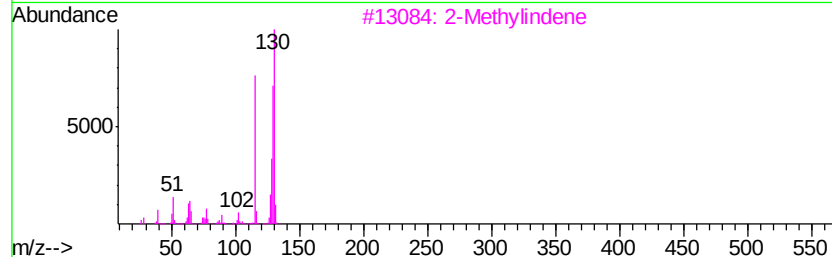
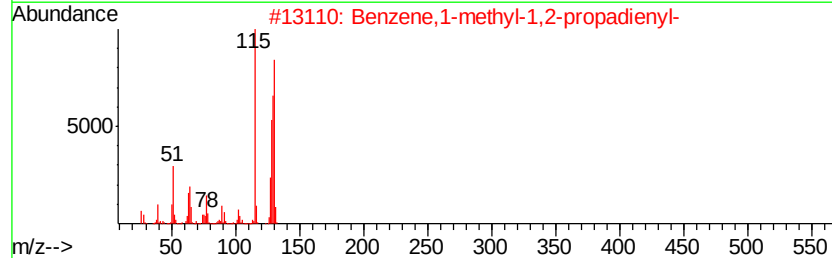
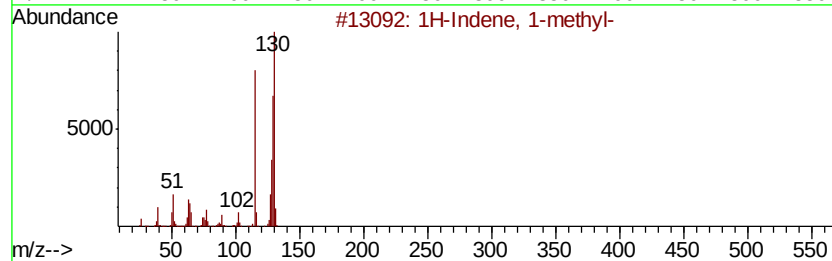
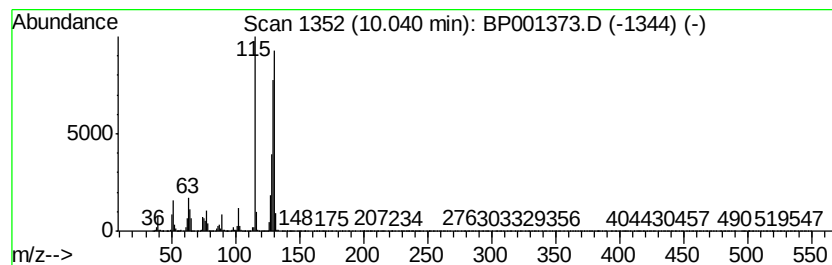
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1H-Indene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.04	2.89 ng	8332090	Naphthalene-d8	10.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 1-methyl-	130 C10H10	000767-59-9 96
2		Benzene,1-methyl-1,2-propadienyl-	130 C10H10	022433-39-2 95
3		2-Methylindene	130 C10H10	002177-47-1 95
4		Cycloprop[alindene, 1,1a,6,6a-te...	130 C10H10	015677-15-3 93
5		Benzene, (1-methyl-2-cyclopropen...	130 C10H10	065051-83-4 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
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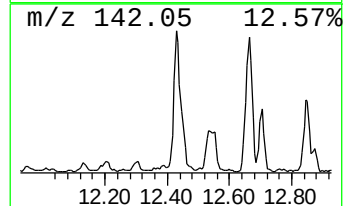
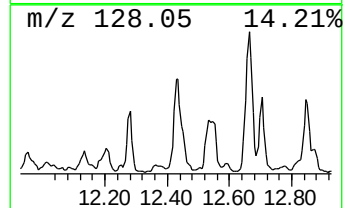
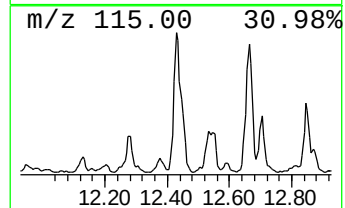
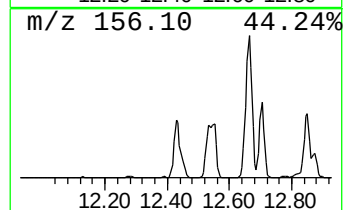
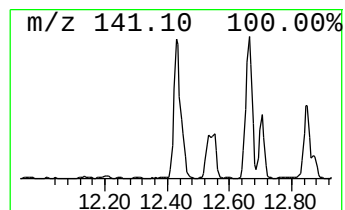
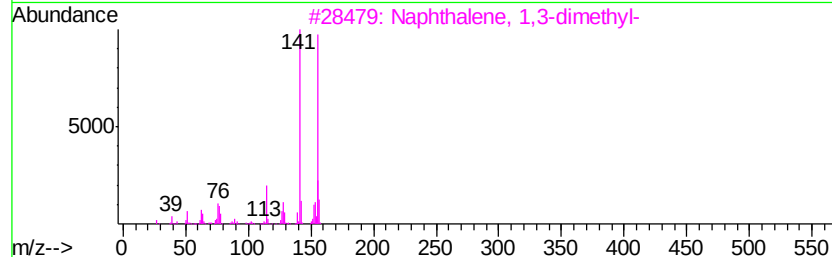
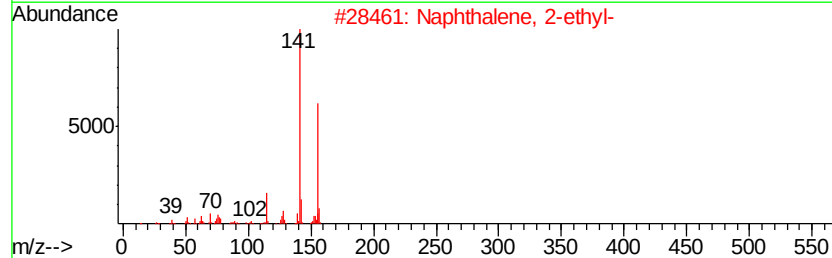
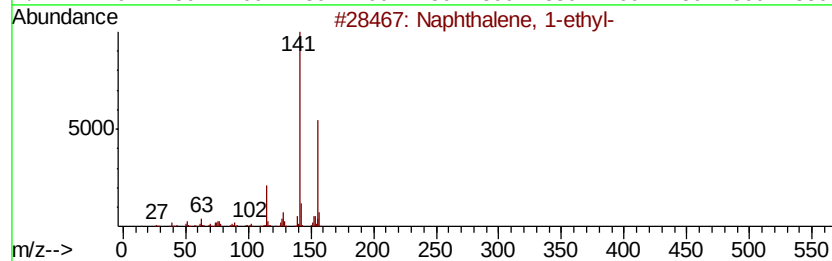
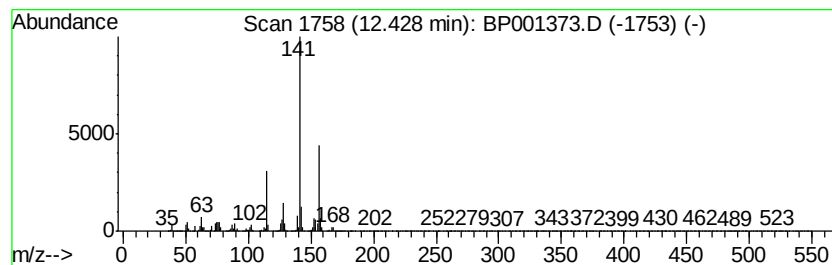
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	29.10 ng	1051850	Acenaphthene-d10	13.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 95
2		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94
3		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 91
4		2-Amino-4-tertbutylthiazole	156 C7H12N2S	074370-93-7 59
5		Methyl phenylphosphinic acid	156 C7H9O2P	004271-13-0 53



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Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

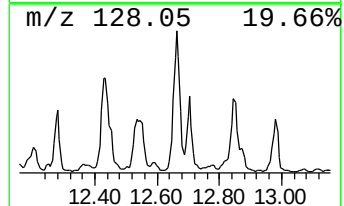
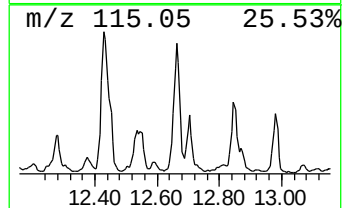
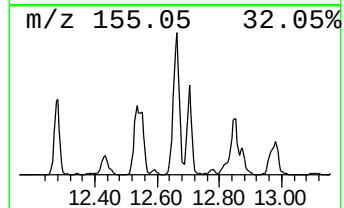
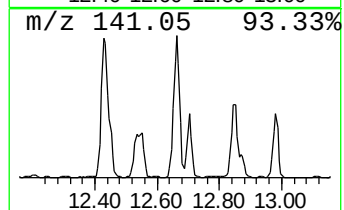
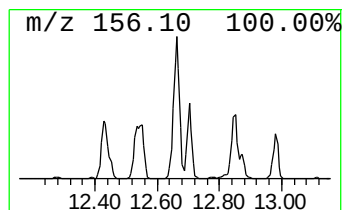
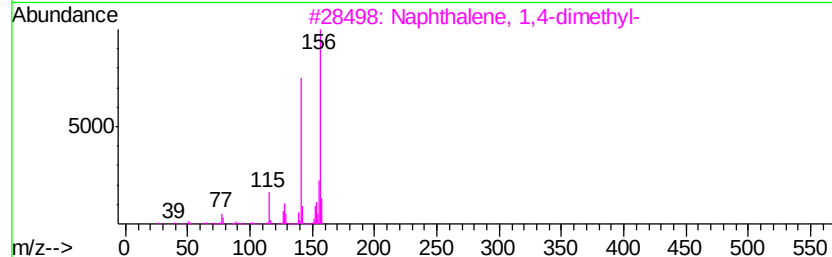
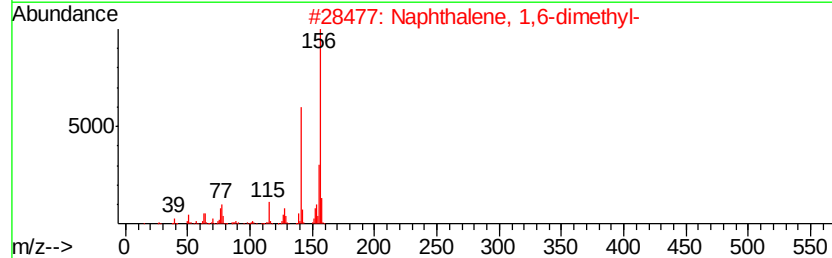
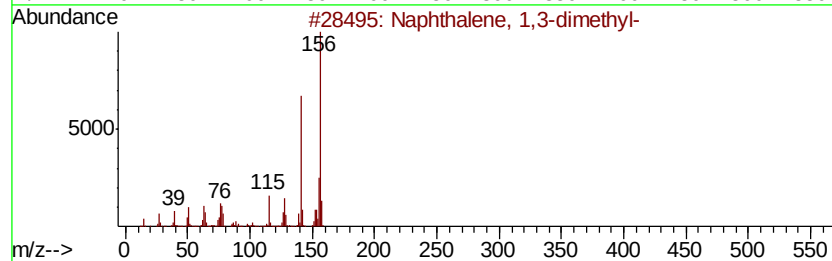
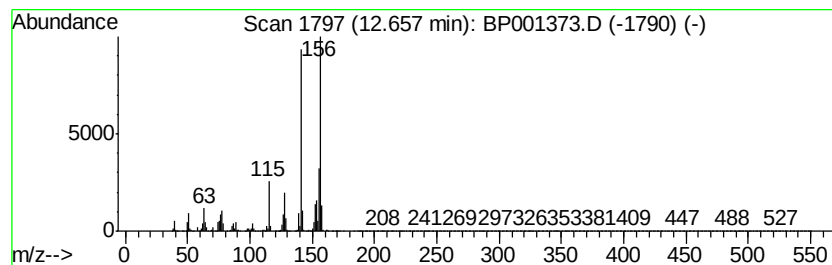
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 1,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.66	35.79 ng	1293590	Acenaphthene-d10		13.20
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	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-14

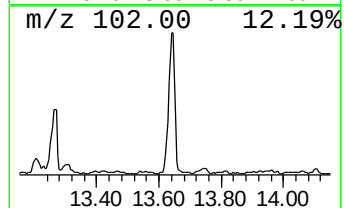
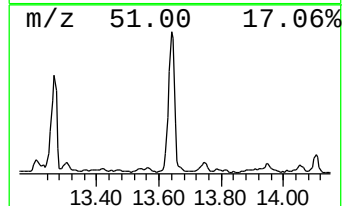
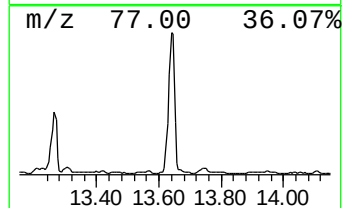
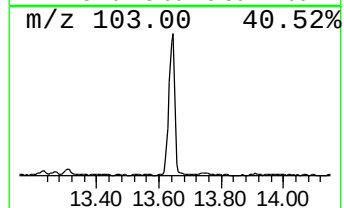
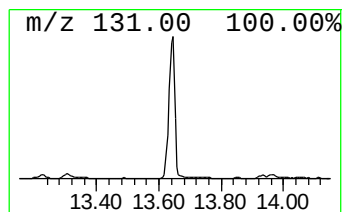
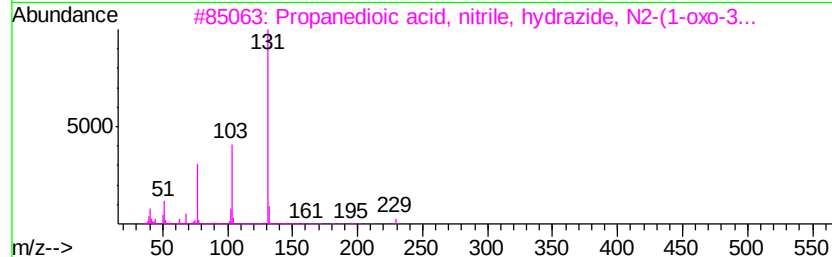
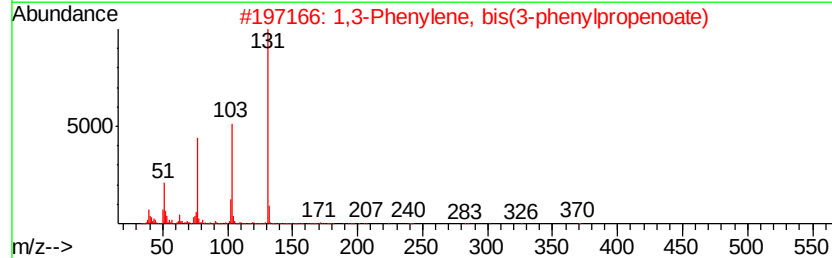
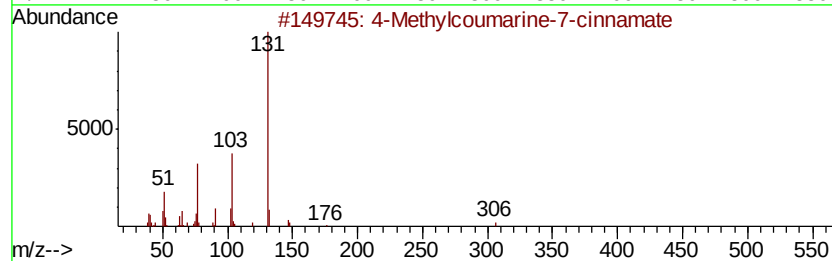
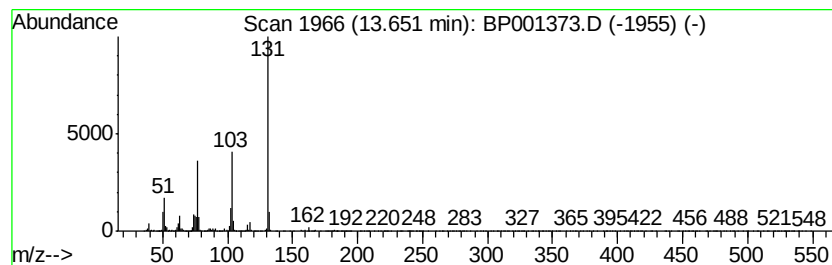
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 4-Methylcoumarine-7-cinnamate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	25.14 ng	908677	Acenaphthene-d10	13.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Methylcoumarine-7-cinnamate	306	C19H14O4	300829-05-4	91
2		1,3-Phenylene, bis(3-phenylprope...	370	C24H18O4	1000259-62-4	83
3		Propanedioic acid, nitrile, hvdr...	229	C12H11N3O2	294878-35-6	72
4		1,3-Bis(cinnamovloxymethyl)adama...	456	C30H32O4	303797-58-2	56
5		1-Penten-3-one, 4-methyl-1-phenyl-	174	C12H14O	003160-32-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP122319\
Data File : BP001373.D
Acq On : 23 Dec 2019 18:08
Operator : JU
Sample : K6372-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-14

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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
Benzene, (1-methy...	7.04	6.8	ng	1323500	1	8.30	3898510	20.0
Benzene, propyl-	7.48	6.0	ng	1159230	1	8.30	3898510	20.0
Benzene, 1-ethyl-...	7.63	5.6	ng	1087880	1	8.30	3898510	20.0
Benzene, 1,2,3-tr...	7.70	3.1	ng	609226	1	8.30	3898510	20.0
Benzene, 1-ethyl-...	7.83	13.8	ng	2688260	1	8.30	3898510	20.0
unknown7.95	7.95	9.3	ng	1803730	1	8.30	3898510	20.0
Indane	8.61	35.5	ng	6913400	1	8.30	3898510	20.0
Benzene, 1,3-diet...	8.72	4.6	ng	891011	1	8.30	3898510	20.0
Benzene, 1-ethyl-...	8.83	5.8	ng	1125800	1	8.30	3898510	20.0
Benzene, 4-ethyl-...	9.06	3.3	ng	639697	1	8.30	3898510	20.0
1H-Indene, 1-methyl-	10.04	2.9	ng	8332090	2	10.39	57602700	20.0
Naphthalene, 1-et...	12.43	29.1	ng	1051850	3	13.20	722824	20.0
Naphthalene, 1,3-...	12.66	35.8	ng	1293590	3	13.20	722824	20.0
4-Methylcoumarine...	13.65	25.1	ng	908677	3	13.20	722824	20.0