

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
 Data File : BM033314.D
 Acq On : 06 Dec 2021 16:11
 Operator : CG/JU
 Sample : M4918-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-39

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_M\Methods\8270-BM112421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM033314.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.119	126	133	147	rBV	21295708	33420923	42.31%	8.341%
2	5.443	350	358	366	rBV	18729505	30460135	38.56%	7.602%
3	5.519	366	371	374	rVV	1038906	1461973	1.85%	0.365%
4	5.584	374	382	391	rVB	27541733	78986209	100.00%	19.712%
5	5.949	434	444	450	rBV	27003289	49885081	63.16%	12.449%
6	6.413	513	523	534	rBV	964485	1493984	1.89%	0.373%
7	6.902	599	606	614	rBV	2482111	3829475	4.85%	0.956%
8	7.013	617	625	630	rVV	9795109	15150992	19.18%	3.781%
9	7.072	630	635	639	rVV	5502408	8201696	10.38%	2.047%
10	7.149	643	648	654	rVB	5090058	7867786	9.96%	1.963%
11	7.313	668	676	688	rBV	6275633	9875941	12.50%	2.465%
12	7.578	712	721	730	rVB	19018875	31799270	40.26%	7.936%
13	7.925	773	780	784	rBV	493810	898005	1.14%	0.224%
14	8.037	792	799	808	rBV	7297465	11869392	15.03%	2.962%
15	8.196	815	826	836	rBV4	356872	899551	1.14%	0.224%
16	8.296	836	843	852	rVV	5493341	9042429	11.45%	2.257%
17	8.407	852	862	866	rVV	556142	983815	1.25%	0.246%
18	8.460	866	871	879	rVV2	1464710	2412518	3.05%	0.602%
19	8.554	879	887	893	rVV	1438049	2428817	3.07%	0.606%
20	8.719	907	915	926	rVB	475732	845805	1.07%	0.211%
21	8.872	934	941	945	rBV	922083	1618470	2.05%	0.404%
22	8.919	945	949	955	rVV	921592	1620691	2.05%	0.404%
23	9.025	955	967	973	rVV2	1898317	4242582	5.37%	1.059%
24	9.096	973	979	987	rVB2	2120143	4566566	5.78%	1.140%
25	9.354	1016	1023	1030	rBV2	532174	951493	1.20%	0.237%
26	9.543	1049	1055	1060	rBV	927914	1428205	1.81%	0.356%
27	9.601	1060	1065	1076	rVB	1290192	2132415	2.70%	0.532%
28	9.966	1122	1127	1140	rVB	1049201	1850532	2.34%	0.462%
29	10.119	1146	1153	1161	rBV2	2600266	4806927	6.09%	1.200%
30	10.354	1187	1193	1199	rVB	557662	948647	1.20%	0.237%
31	10.719	1244	1255	1258	rBV2	535368	1345269	1.70%	0.336%
32	10.778	1258	1265	1271	rVV	9543857	16579335	20.99%	4.138%
33	12.107	1484	1491	1498	rBV	970992	1737010	2.20%	0.433%
34	12.378	1525	1537	1543	rBV	5215060	8705240	11.02%	2.172%
35	12.595	1564	1574	1580	rBV	4896106	8489102	10.75%	2.119%
36	13.172	1662	1672	1679	rBV	3982480	5944688	7.53%	1.484%

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.378	1700	1707	1713	rBV2	498783	1030583	1.30%	0.257%
38	13.572	1735	1740	1754	rVB2	683838	1650333	2.09%	0.412%
39	13.725	1754	1766	1773	rBV4	1499272	4439612	5.62%	1.108%
40	13.866	1784	1790	1795	rVV	1737494	3271619	4.14%	0.816%
41	13.925	1795	1800	1811	rVB	1195284	2091682	2.65%	0.522%
42	14.107	1825	1831	1834	rBV	691883	1048127	1.33%	0.262%
43	14.548	1896	1906	1912	rBV2	1013945	2077536	2.63%	0.518%
44	16.036	2153	2159	2165	rBV	2862439	4293444	5.44%	1.071%
45	17.283	2366	2371	2376	rBV	1131515	1693554	2.14%	0.423%
46	19.895	2810	2815	2822	rBV	5304091	6712710	8.50%	1.675%
47	21.448	3074	3079	3085	rBV	1415131	1776529	2.25%	0.443%
48	23.777	3469	3475	3484	rVB	912412	1839051	2.33%	0.459%

Sum of corrected areas: 400705749

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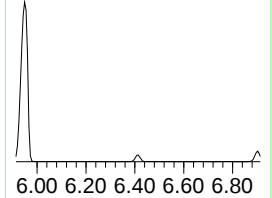
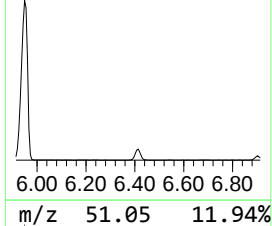
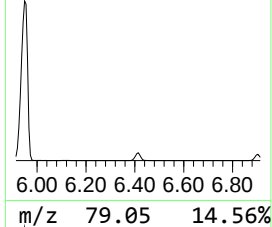
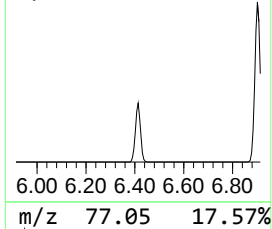
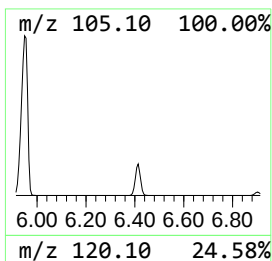
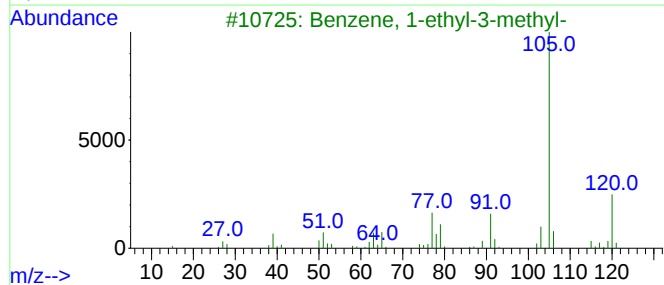
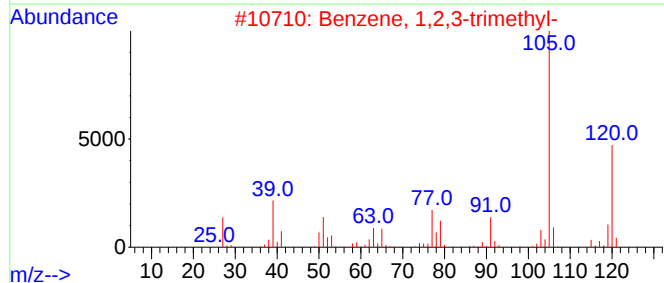
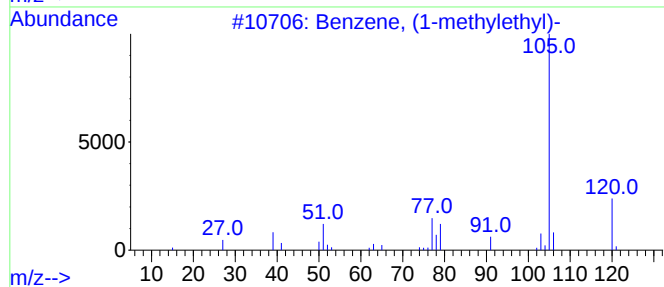
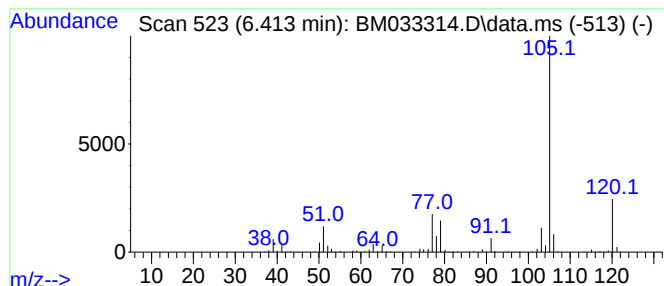
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, (1-methylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.413	33.27 ng	1493980	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	95
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	83
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80



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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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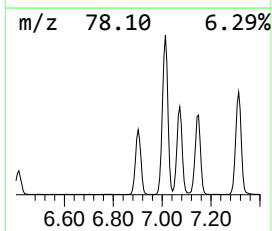
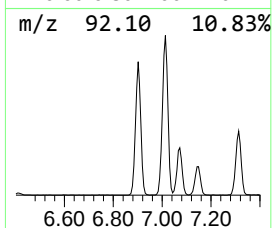
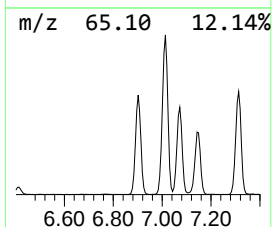
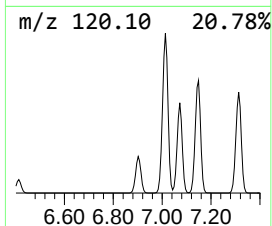
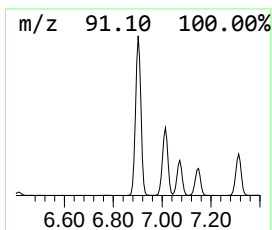
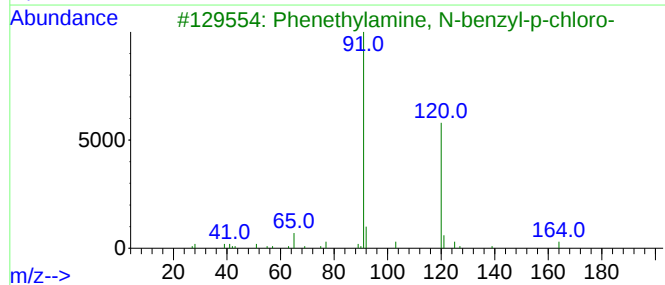
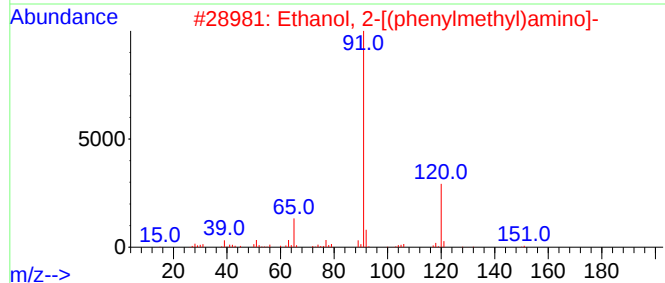
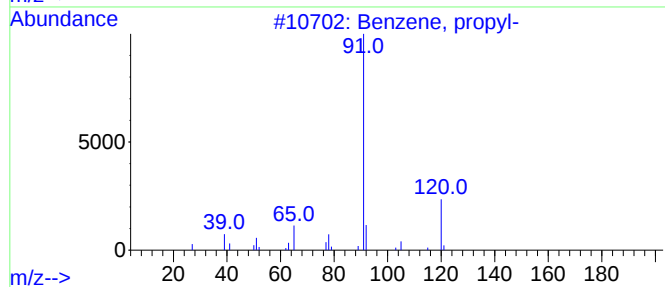
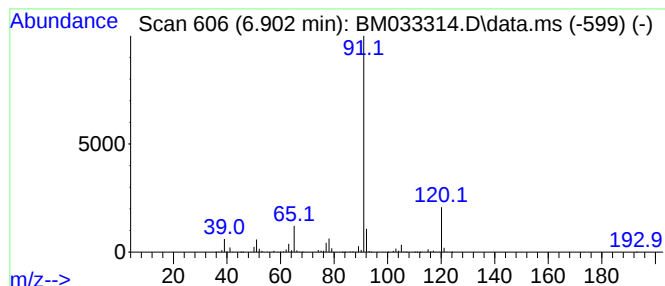
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.902	85.29 ng	3829480	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	78
3			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78
4			1,2-Ethanediamine, N,N'-bis(phen...)	240	C16H20N2	000140-28-3	72
5			1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	53



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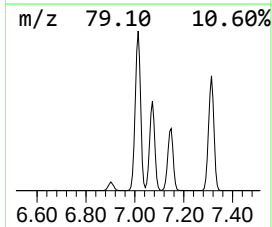
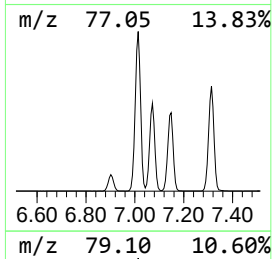
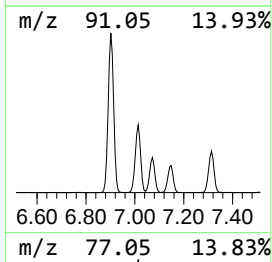
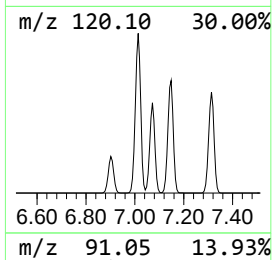
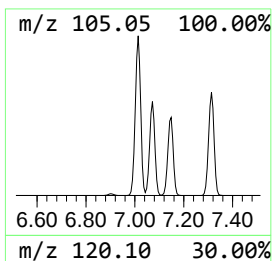
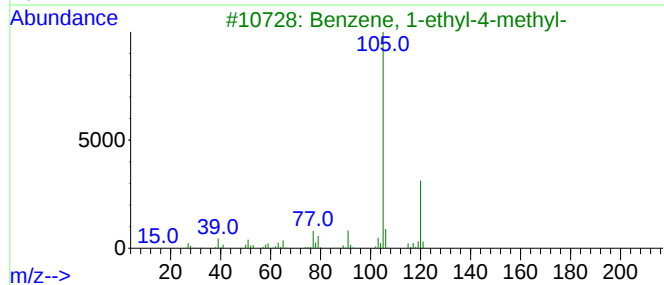
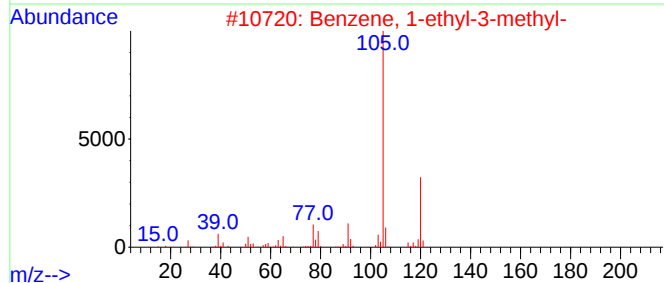
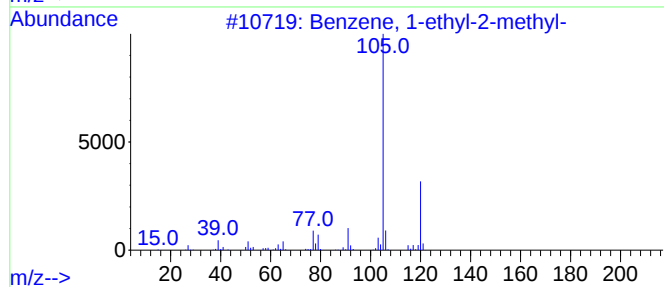
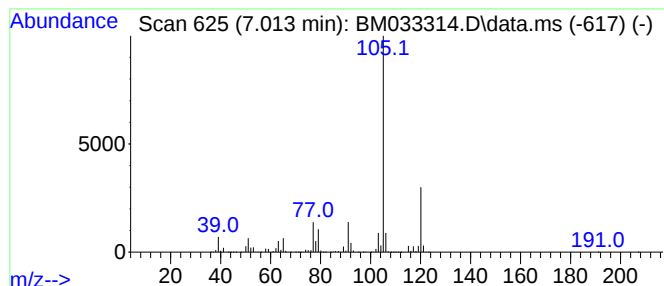
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.013	337.44 ng	15151000	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Instrument :
BNA_M
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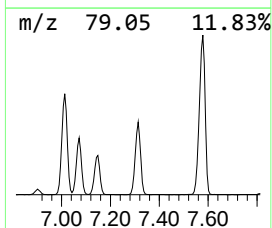
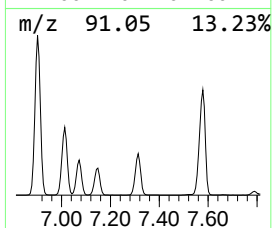
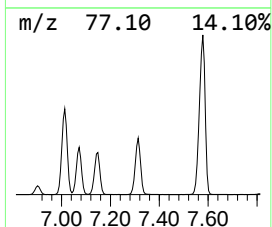
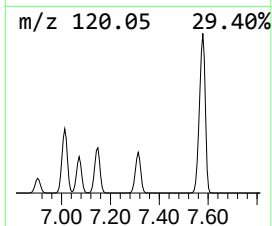
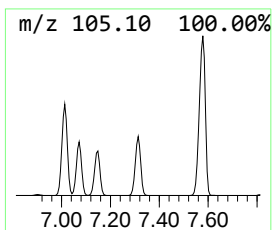
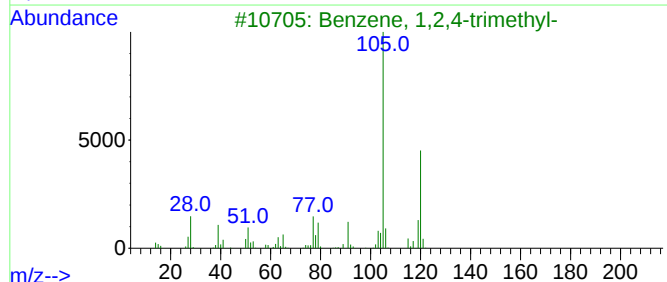
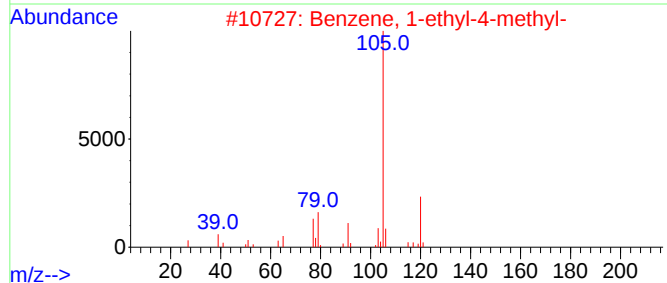
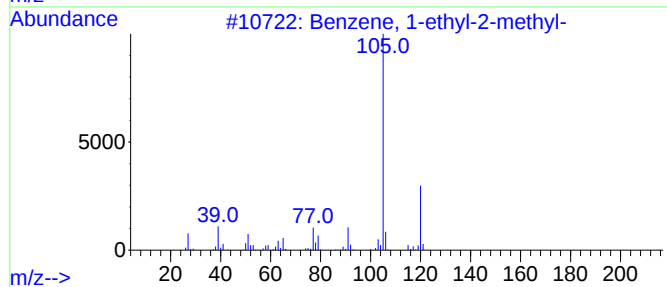
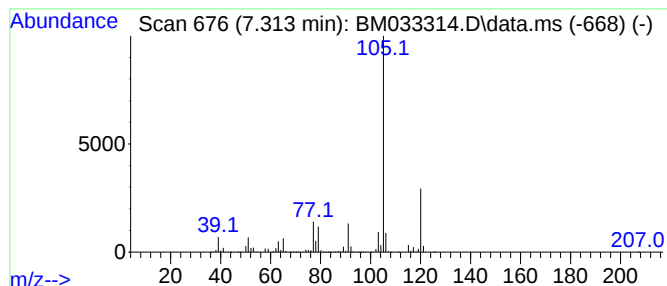
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.313	219.95 ng	9875940	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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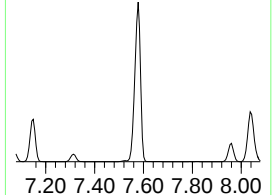
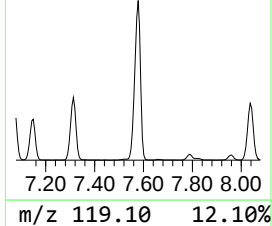
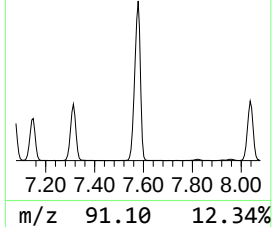
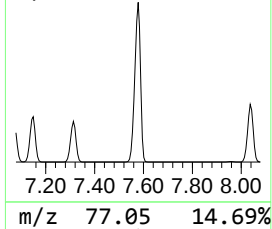
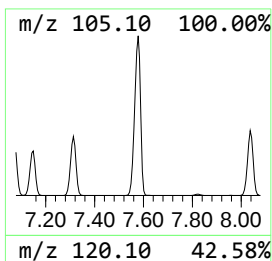
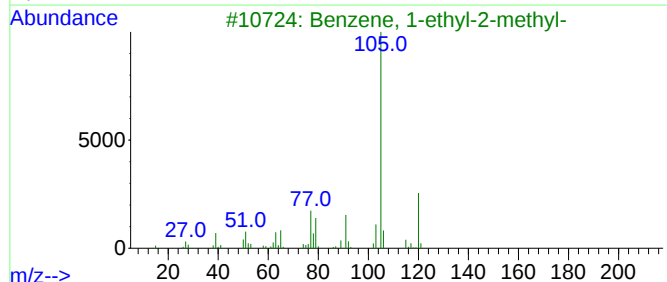
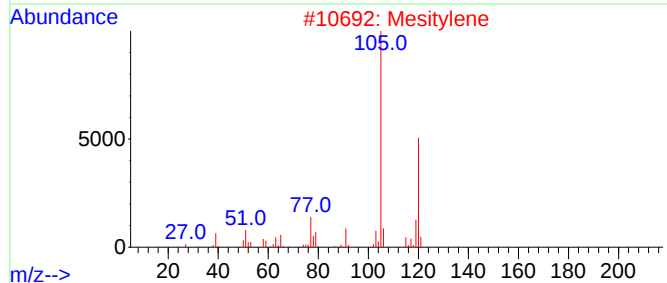
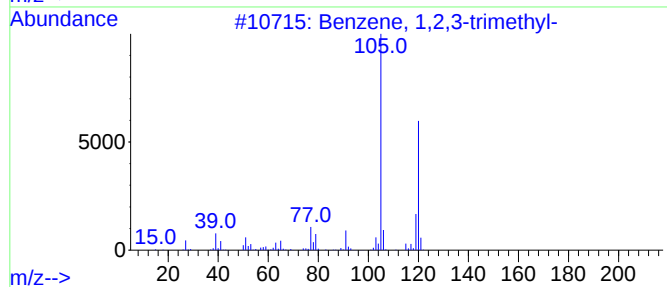
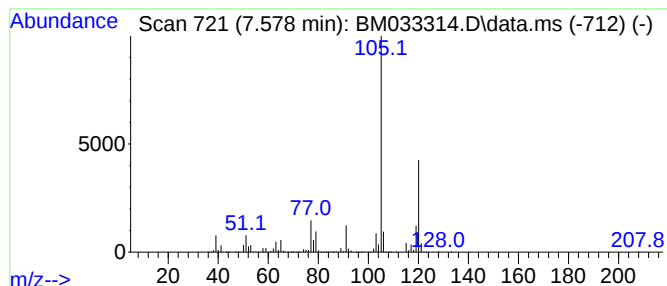
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.578	708.22 ng	31799300	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	97
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
Operator : CG/JU
Sample : M4918-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-39

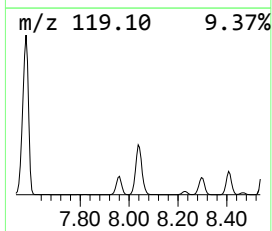
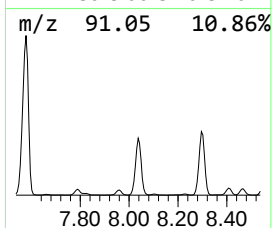
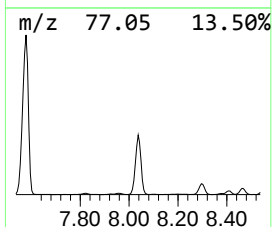
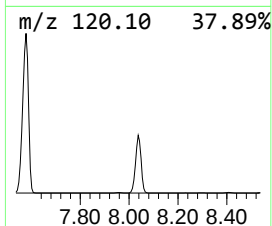
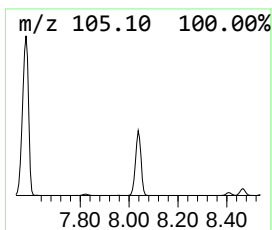
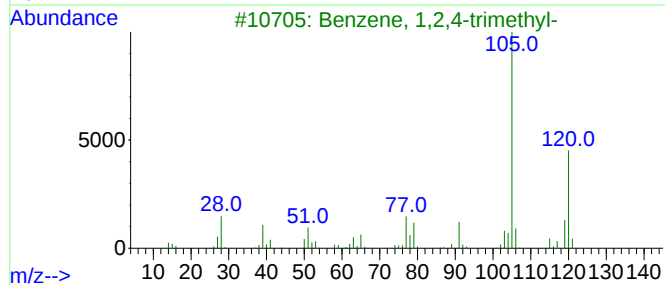
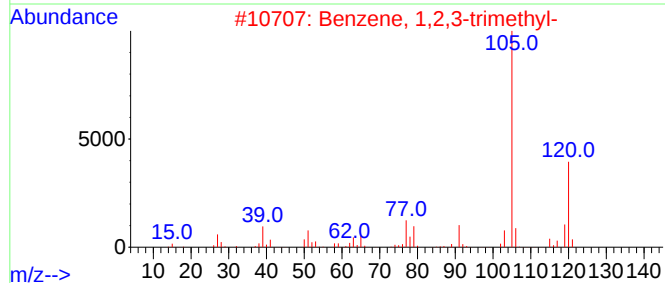
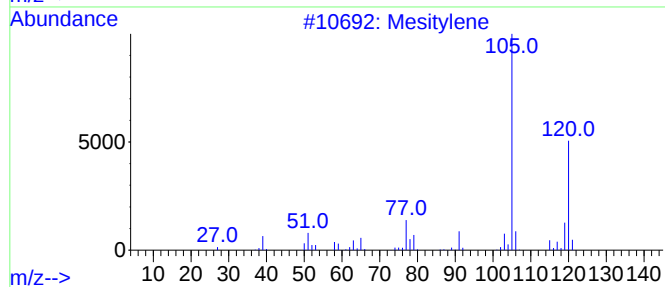
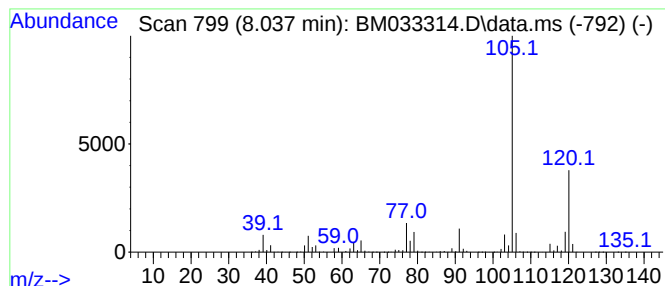
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.037	264.35 ng	11869400	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
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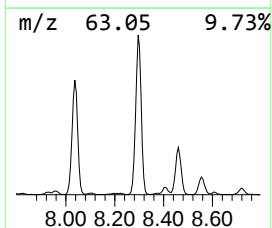
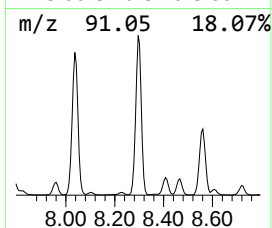
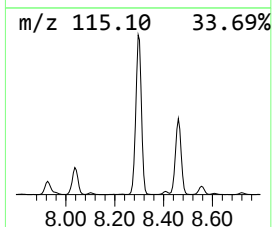
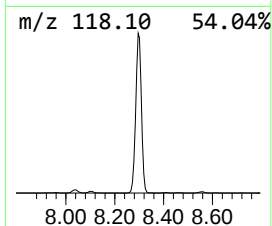
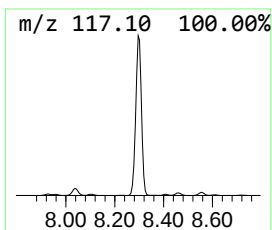
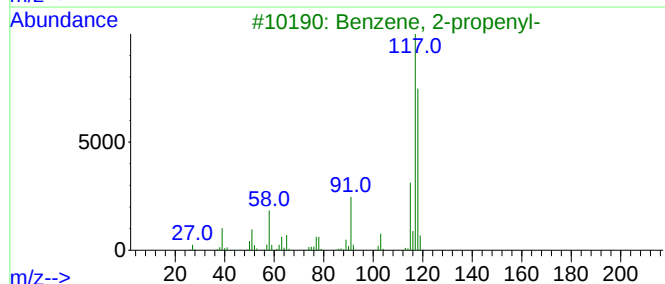
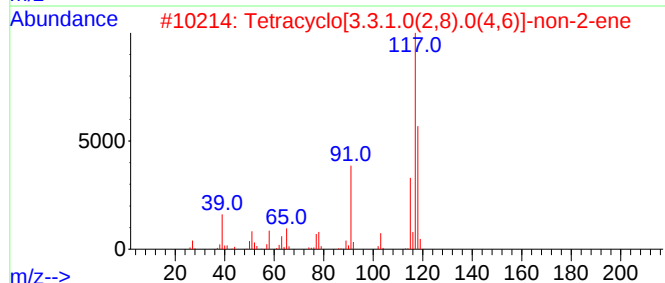
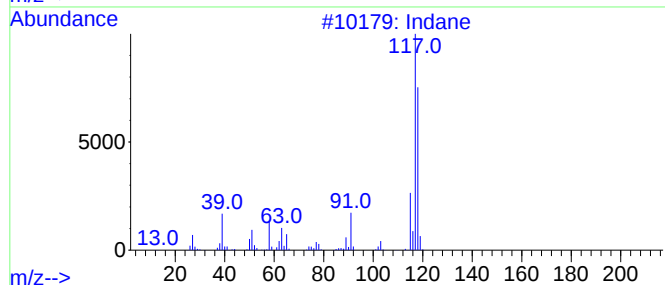
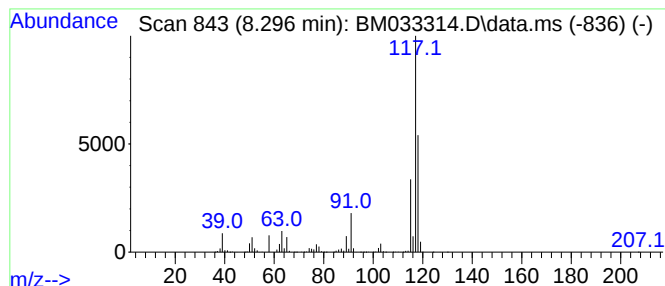
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.296	201.39 ng	9042430	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
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Sample : M4918-04
Misc :
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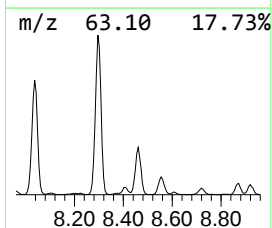
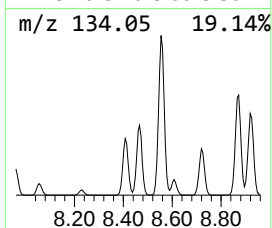
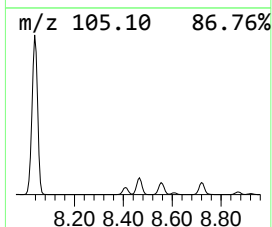
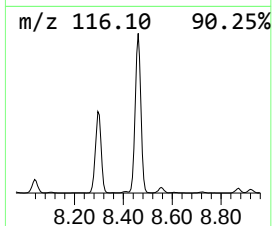
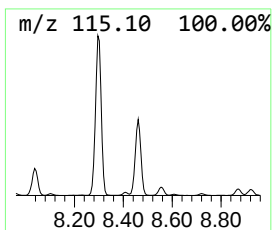
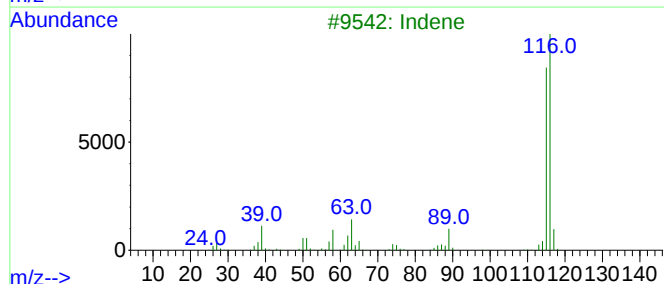
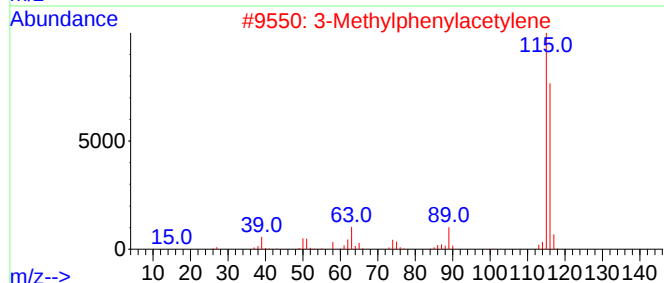
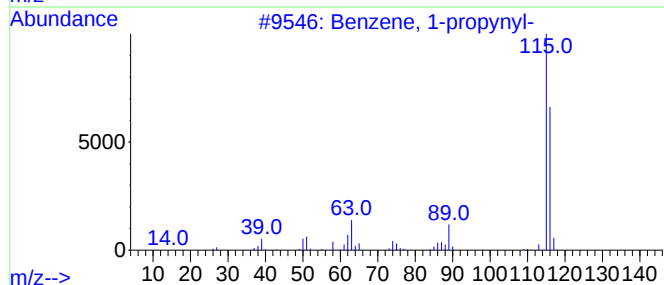
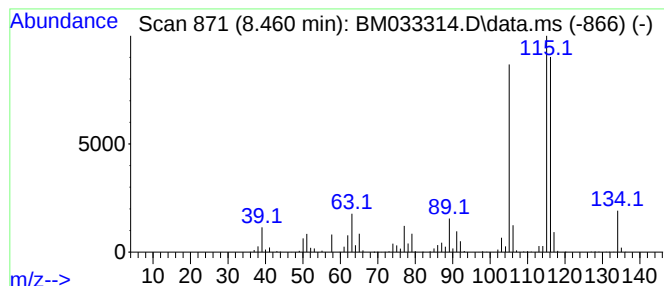
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1-propynyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.460	53.73 ng	2412520	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
2			3-Methylphenylacetylene	116	C9H8	000766-82-5	93
3			Indene	116	C9H8	000095-13-6	89
4			2-Methylphenylacetylene	116	C9H8	000766-47-2	68
5			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
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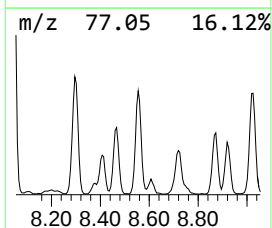
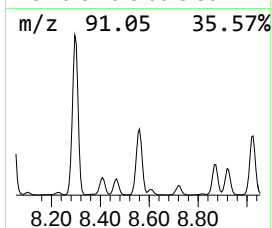
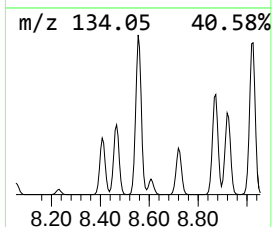
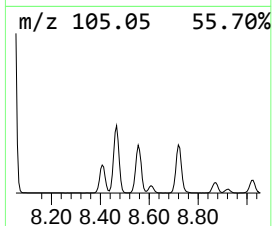
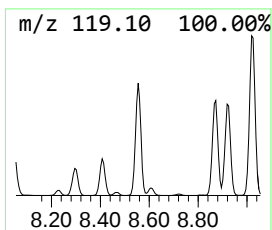
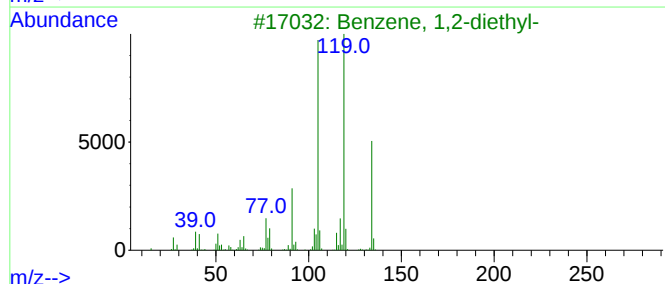
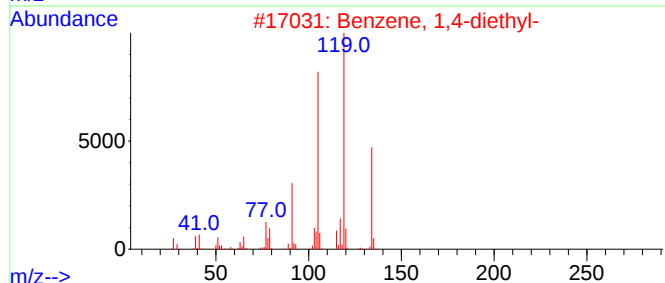
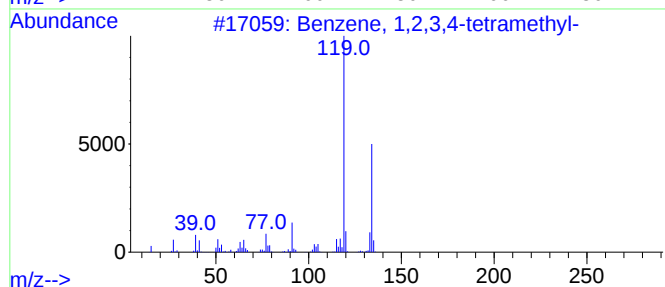
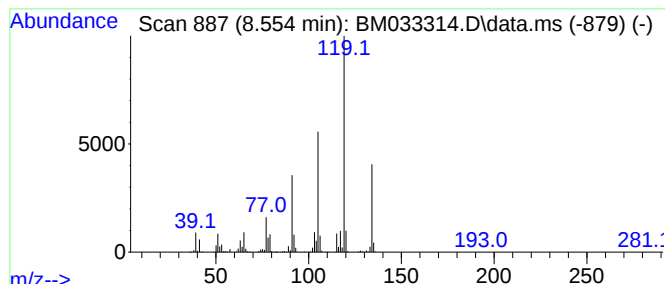
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.554	54.09 ng	2428820	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
4			o-Cymene	134	C10H14	000527-84-4	93
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
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ClientSampleId :
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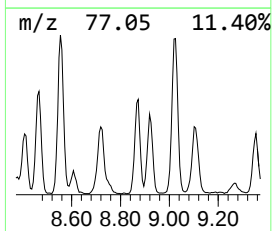
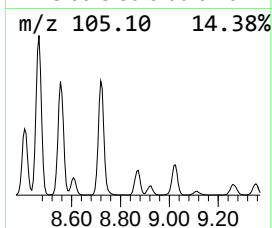
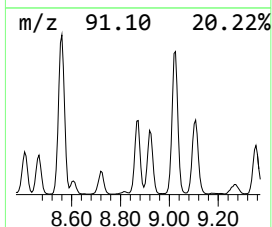
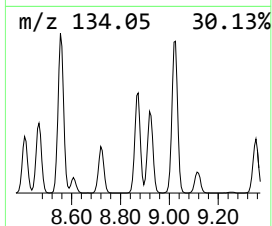
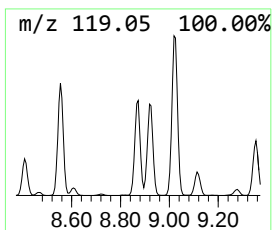
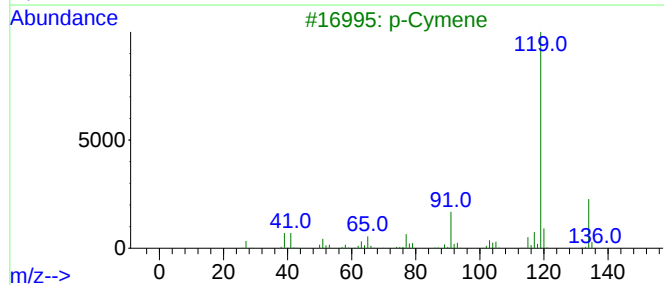
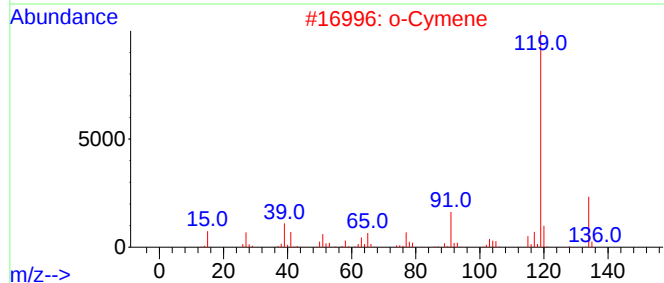
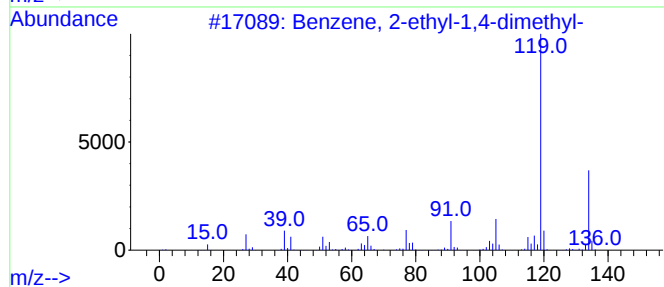
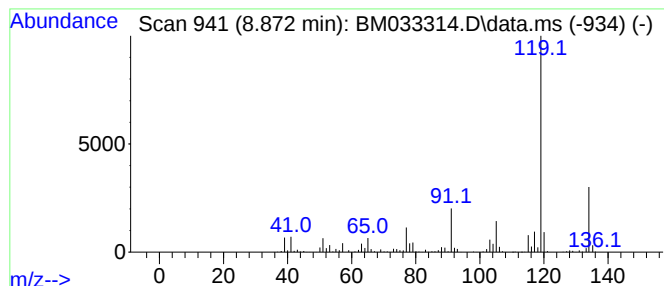
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.872	36.05 ng	1618470	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



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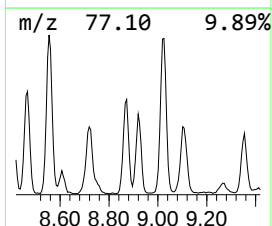
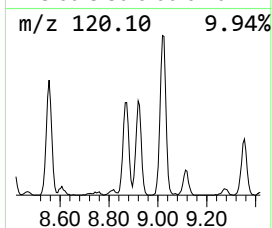
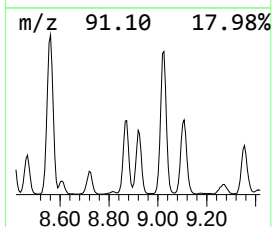
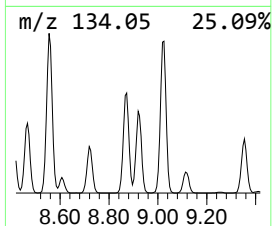
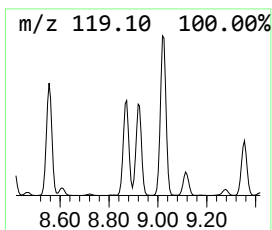
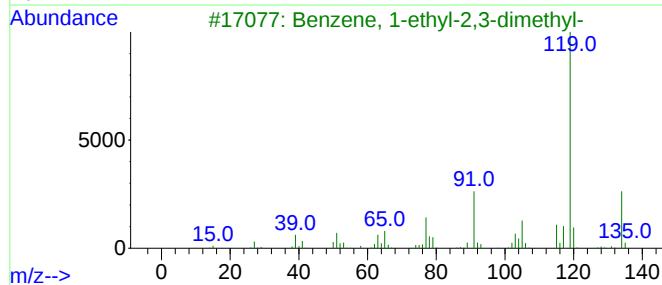
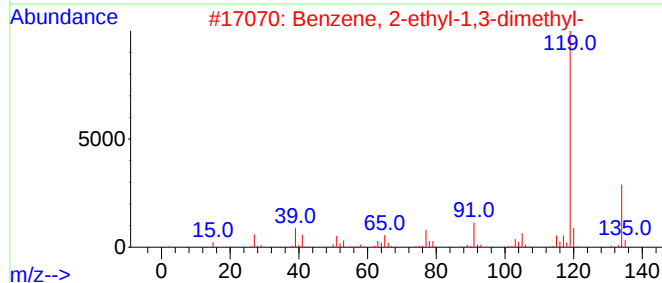
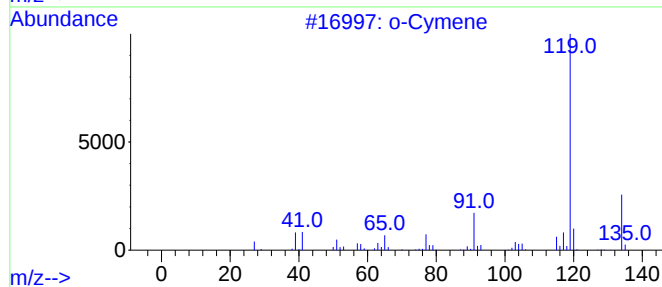
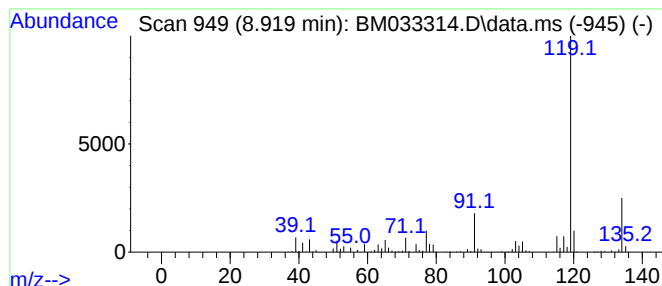
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 o-Cymene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.919	36.10 ng	1620690	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
4			p-Cymene	134	C10H14	000099-87-6	94
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
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Misc :
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ClientSampleId :
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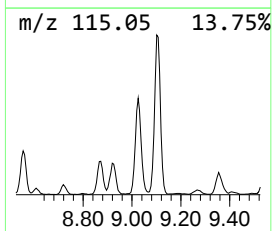
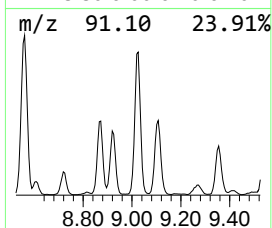
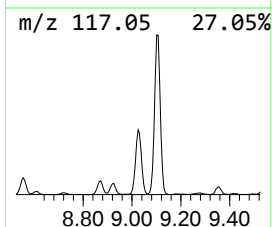
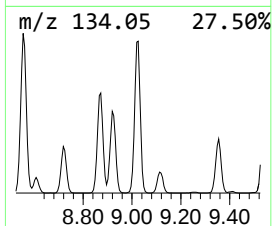
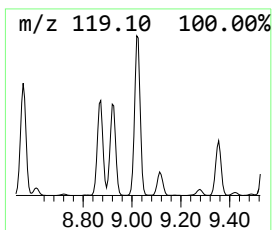
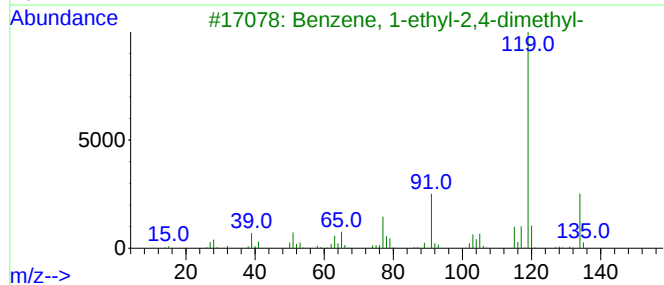
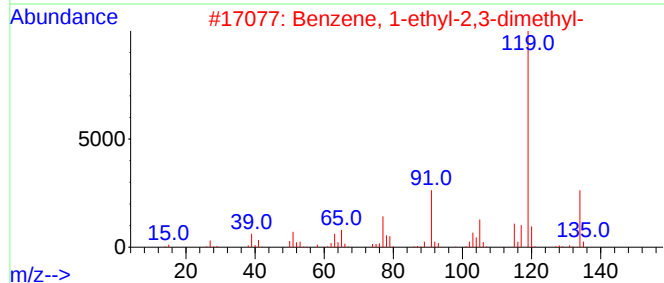
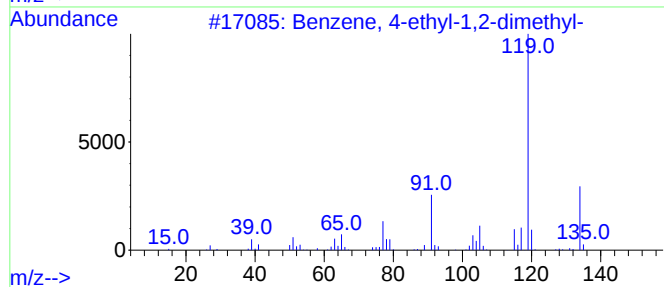
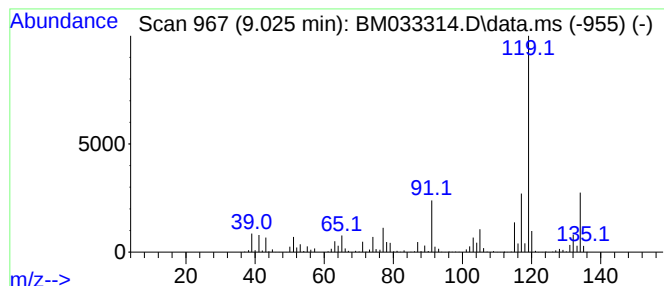
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.025	94.49 ng	4242580	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
Operator : CG/JU
Sample : M4918-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MW-39

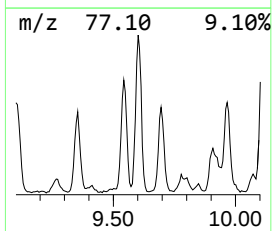
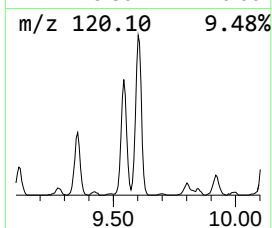
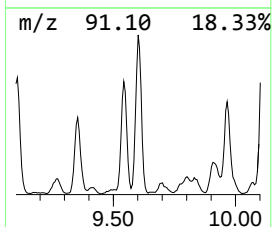
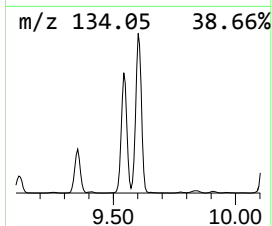
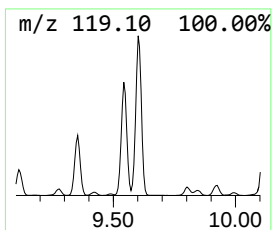
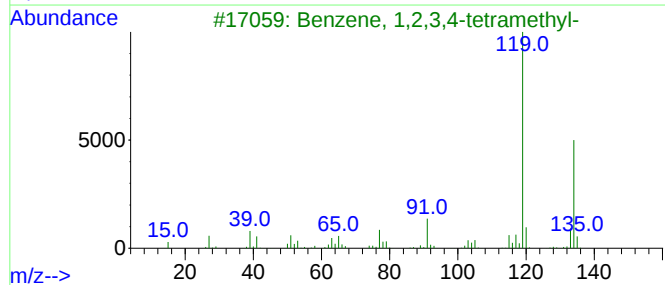
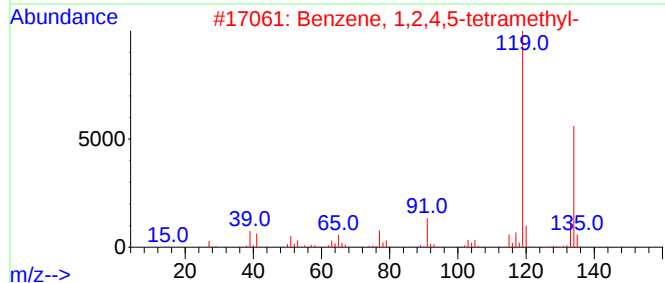
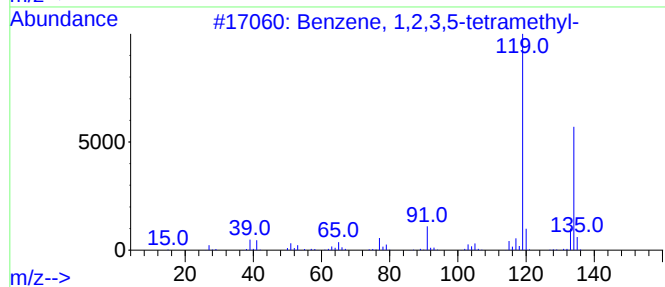
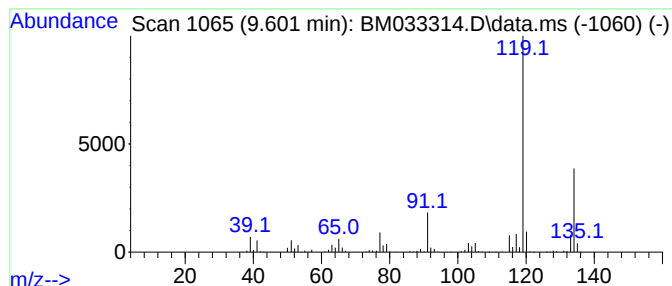
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.601	31.70 ng	2132420	Naphthalene-d8	10.725

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
Operator : CG/JU
Sample : M4918-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MW-39

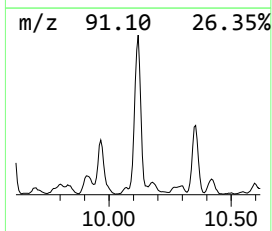
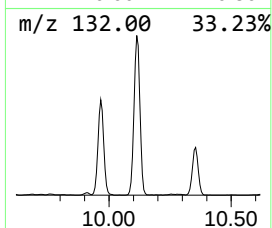
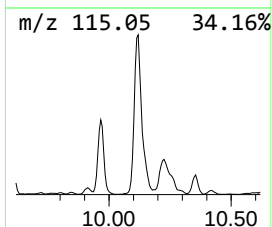
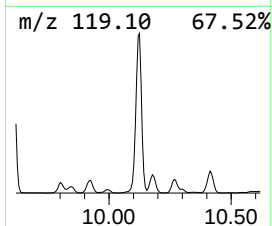
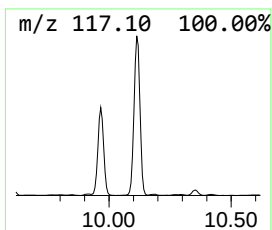
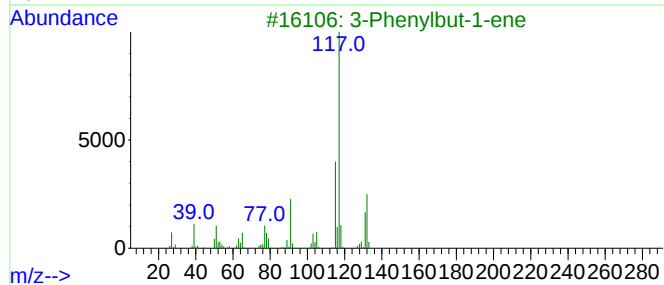
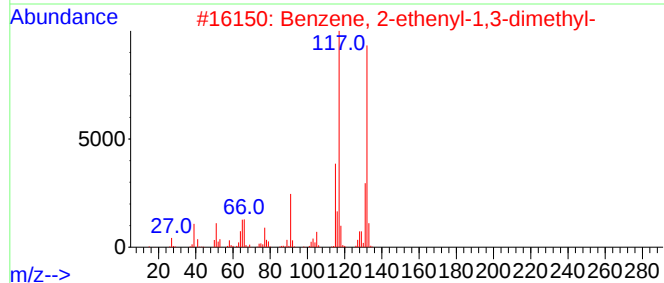
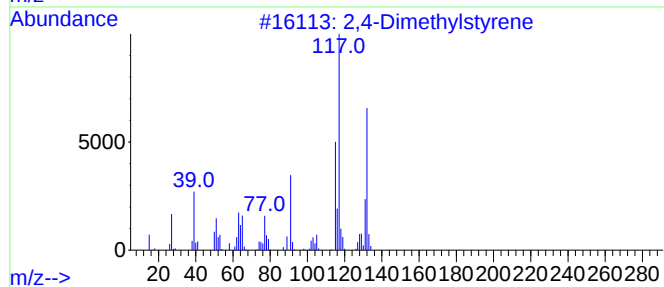
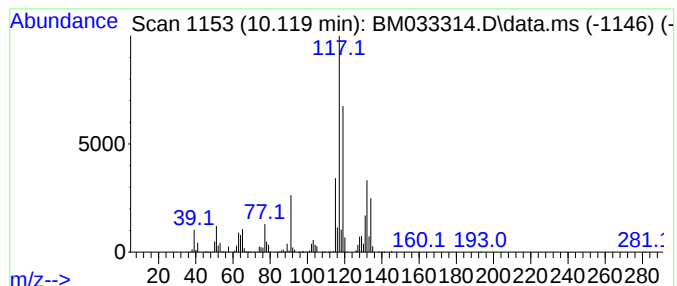
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 2,4-Dimethylstyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.119	71.46 ng	4806930	Naphthalene-d8	10.725

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
Operator : CG/JU
Sample : M4918-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MW-39

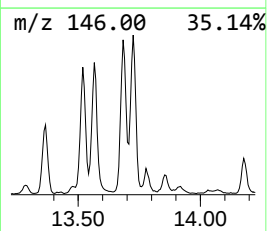
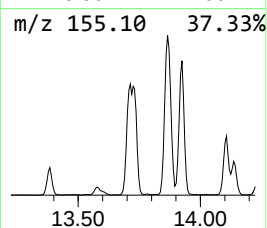
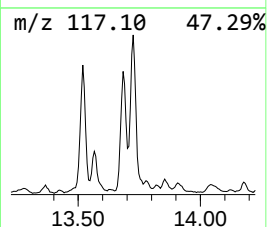
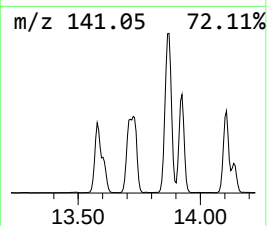
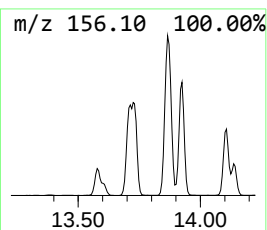
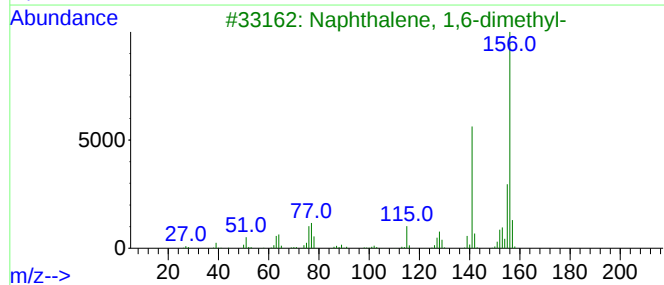
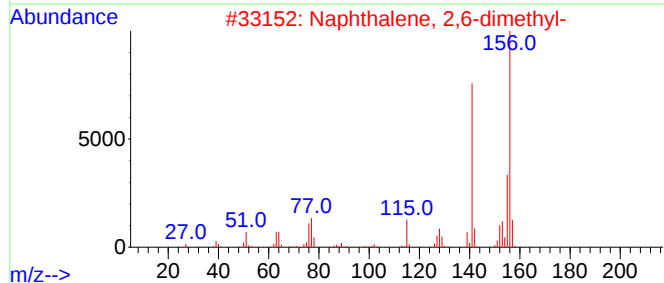
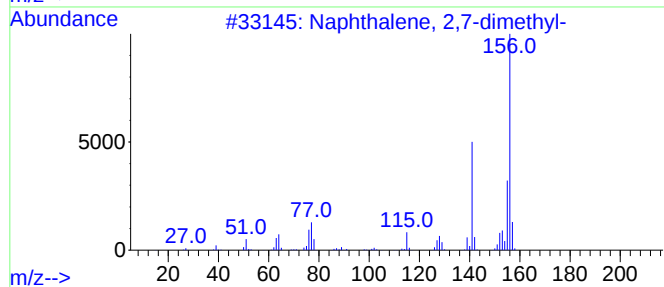
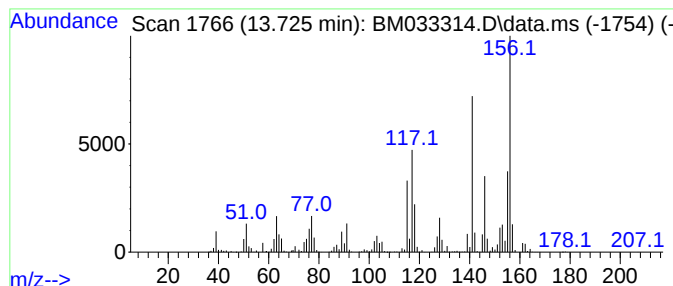
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Naphthalene, 2,7-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.725	42.74 ng	4439610	Acenaphthene-d10	14.548

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
Data File : BM033314.D
Acq On : 06 Dec 2021 16:11
Operator : CG/JU
Sample : M4918-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

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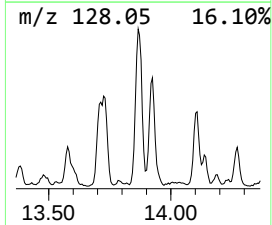
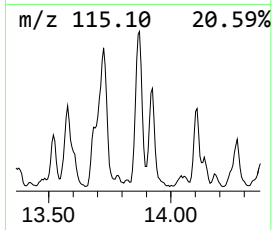
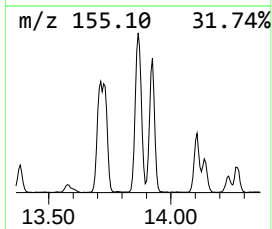
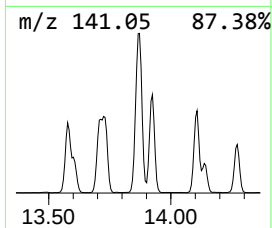
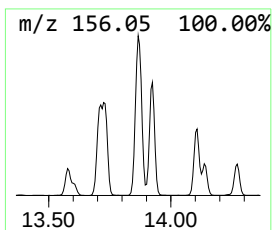
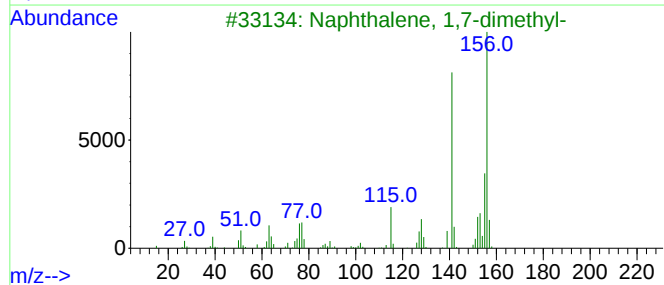
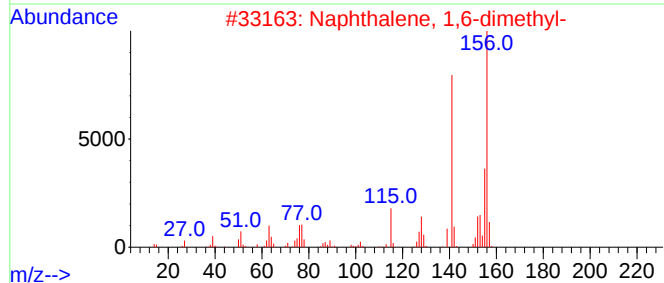
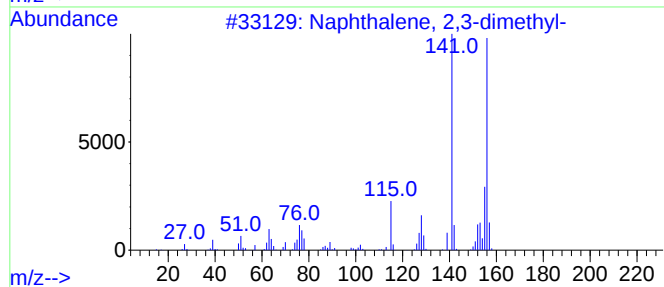
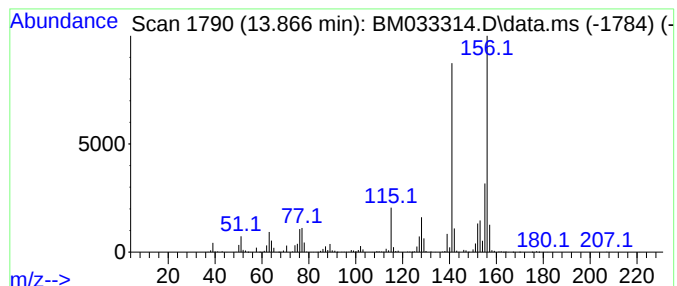
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Naphthalene, 2,3-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.866	31.50 ng	3271620	Acenaphthene-d10	14.548

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
 Data File : BM033314.D
 Acq On : 06 Dec 2021 16:11
 Operator : CG/JU
 Sample : M4918-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.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, propyl-	6.902	85.3	ng	3829480	1	7.925	898005	20.0
Benzene, 1-ethy...	7.013	337.4	ng	15151000	1	7.925	898005	20.0
Benzene, 1-ethy...	7.313	219.9	ng	9875940	1	7.925	898005	20.0
Benzene, 1,2,3-...	7.578	708.2	ng	31799300	1	7.925	898005	20.0
Mesitylene	8.037	264.4	ng	11869400	1	7.925	898005	20.0
Indane	8.296	201.4	ng	9042430	1	7.925	898005	20.0
Benzene, 1-prop...	8.460	53.7	ng	2412520	1	7.925	898005	20.0
Benzene, 1,2,3,...	8.554	54.1	ng	2428820	1	7.925	898005	20.0
Benzene, 2-ethy...	8.872	36.0	ng	1618470	1	7.925	898005	20.0
o-Cymene	8.919	36.1	ng	1620690	1	7.925	898005	20.0
Benzene, 4-ethy...	9.025	94.5	ng	4242580	1	7.925	898005	20.0
Benzene, 1,2,3,...	9.601	31.7	ng	2132420	2	10.725	1345270	20.0
2,4-Dimethylsty...	10.119	71.5	ng	4806930	2	10.725	1345270	20.0
Naphthalene, 2,...	13.725	42.7	ng	4439610	3	14.548	2077540	20.0
Naphthalene, 2,...	13.866	31.5	ng	3271620	3	14.548	2077540	20.0