

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
 Data File : BM036198.D
 Acq On : 10 Aug 2022 18:46
 Operator : CG/JU
 Sample : N3971-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-14R

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-BM080122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM036198.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.981	147	152	165	rVB	2182332	2864903	6.12%	1.363%
2	4.281	199	203	209	rVV	1231403	1724902	3.69%	0.821%
3	4.934	307	314	321	rBV3	475967	787242	1.68%	0.375%
4	5.322	373	380	390	rBV	8559396	11976483	25.59%	5.698%
5	5.416	390	396	399	rVV	2900709	4316961	9.22%	2.054%
6	5.463	399	404	419	rVB	21217227	46803252	100.00%	22.266%
7	5.834	459	467	478	rBV	14386290	21932579	46.86%	10.434%
8	6.310	535	548	556	rBV	557748	864976	1.85%	0.411%
9	6.798	626	631	637	rVB	1253027	1824722	3.90%	0.868%
10	6.910	637	650	656	rVV	5652972	8536626	18.24%	4.061%
11	6.969	656	660	665	rVV	2368161	3413492	7.29%	1.624%
12	7.045	665	673	685	rVB2	3249985	6605455	14.11%	3.142%
13	7.216	694	702	712	rBV	3469241	5269924	11.26%	2.507%
14	7.475	735	746	754	rBV	10058558	15019202	32.09%	7.145%
15	7.834	798	807	815	rBV	1619494	2640934	5.64%	1.256%
16	7.939	820	825	833	rVB	3861897	5992076	12.80%	2.851%
17	8.198	863	869	880	rBV	1866598	2982910	6.37%	1.419%
18	8.469	905	915	920	rBV2	353617	650817	1.39%	0.310%
19	8.933	987	994	999	rBV	434040	684874	1.46%	0.326%
20	9.004	999	1006	1015	rVB	2769167	4573000	9.77%	2.176%
21	10.028	1176	1180	1190	rVB2	567324	1062616	2.27%	0.506%
22	10.633	1276	1283	1287	rBV	2019480	3419161	7.31%	1.627%
23	10.686	1287	1292	1299	rVB	3034047	4881650	10.43%	2.322%
24	12.298	1555	1566	1572	rBV	553031	897399	1.92%	0.427%
25	12.521	1591	1604	1617	rVB2	417382	798664	1.71%	0.380%
26	13.104	1693	1703	1712	rBV	6397454	9359294	20.00%	4.452%
27	13.333	1736	1742	1755	rVB	1487341	2089141	4.46%	0.994%
28	14.474	1928	1936	1944	rBV	2888486	4222897	9.02%	2.009%
29	15.962	2178	2189	2199	rBV	4781058	6748330	14.42%	3.210%
30	17.215	2395	2402	2409	rBV	3184798	4409865	9.42%	2.098%
31	19.839	2843	2848	2856	rBV	10003261	12396534	26.49%	5.897%
32	21.397	3108	3113	3121	rBV	4082744	5009543	10.70%	2.383%
33	23.744	3505	3512	3523	rVB2	2809146	5442891	11.63%	2.589%

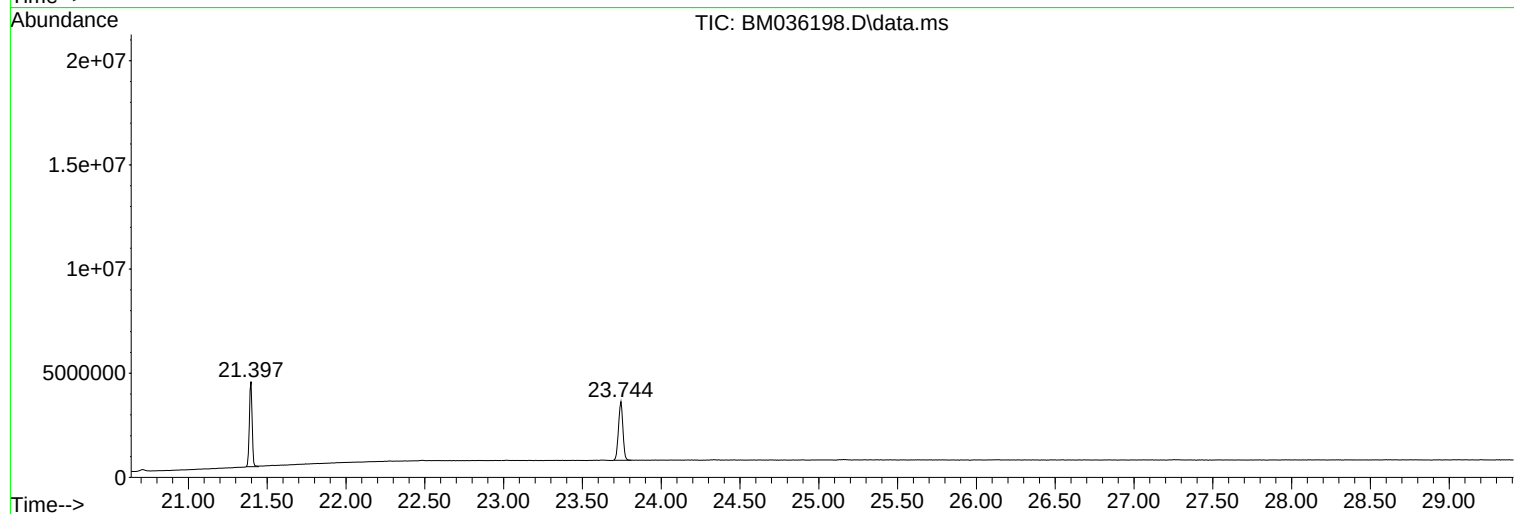
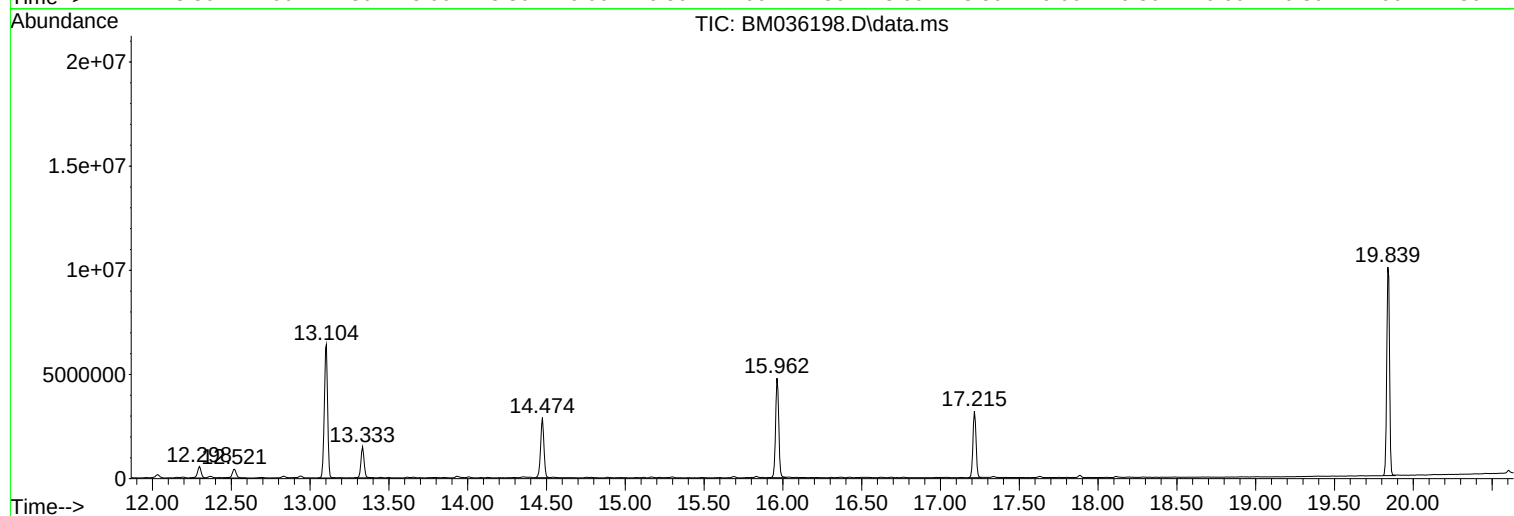
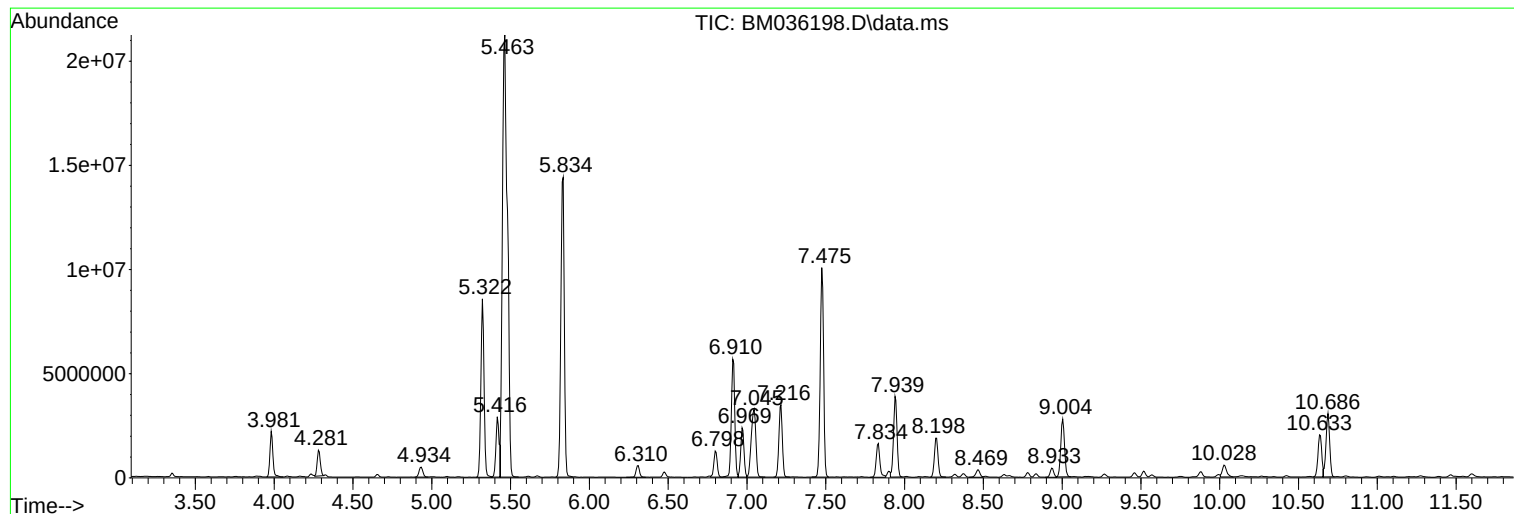
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TIC Library : C:\Database\NIST20.L
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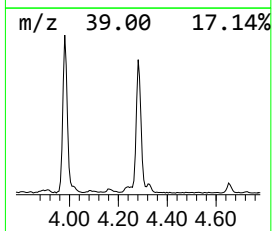
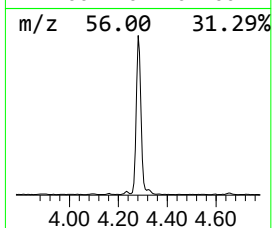
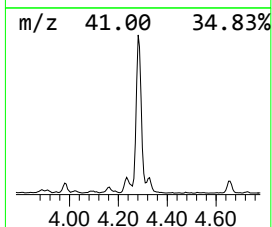
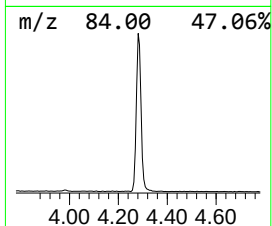
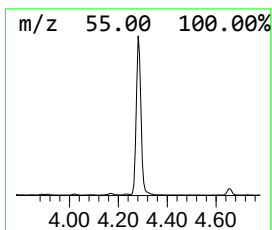
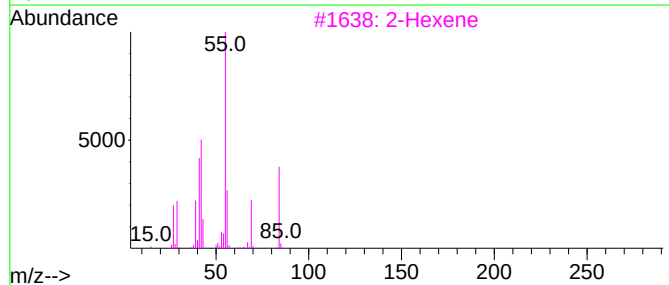
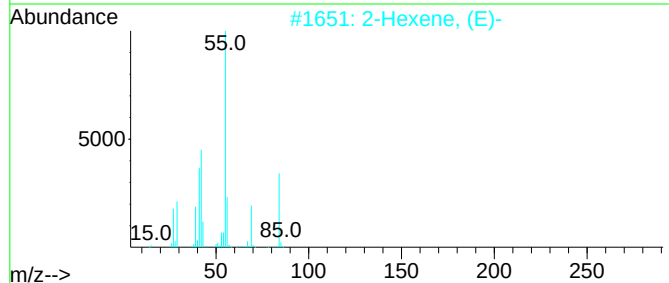
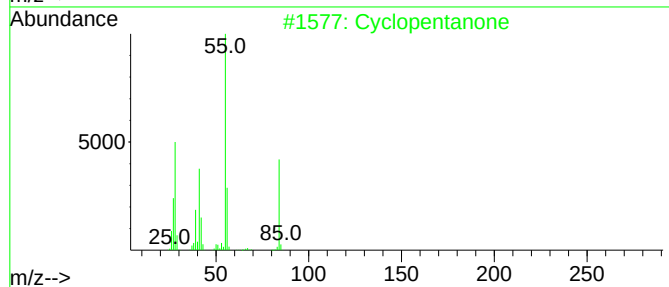
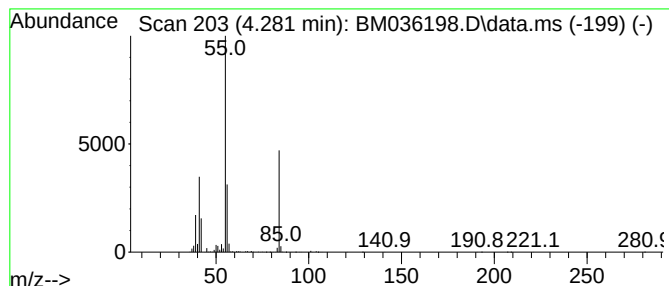
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Cyclopentanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.281	13.06 ng	1724900	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone	84	C5H8O	000120-92-3	91
2			2-Hexene, (E)-	84	C6H12	004050-45-7	59
3			2-Hexene	84	C6H12	000592-43-8	59
4			2(5H)-Furanone	84	C4H4O2	000497-23-4	49
5			3-Hexene	84	C6H12	000592-47-2	45



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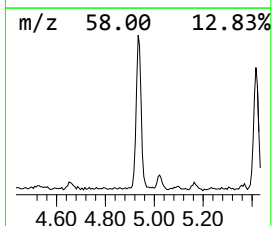
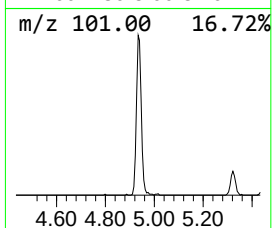
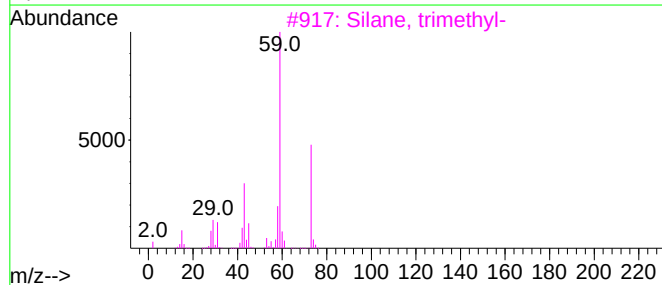
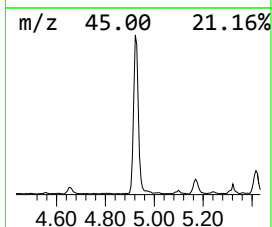
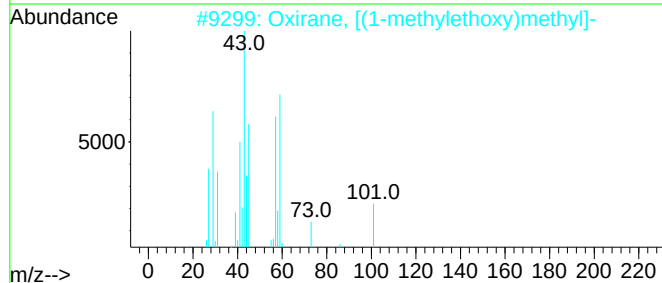
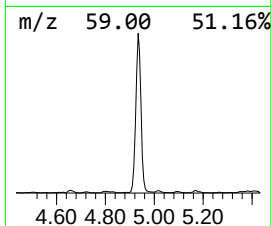
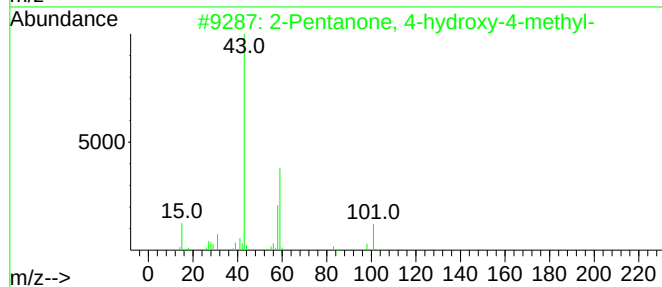
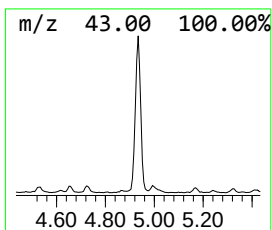
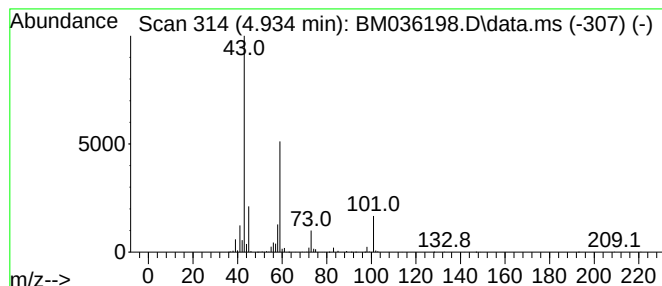
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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	5.96 ng	787242	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
2			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	45
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	43
4			2-(2-methoxyethoxy)propanoic aci...	190	C8H14O5	1000506-61-7	36
5			1,3,4-Oxadiazole, 2-(acetyloxy)-...	172	C7H12N2O3	066398-57-0	35



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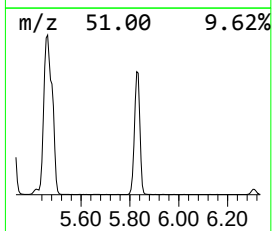
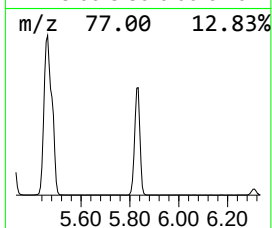
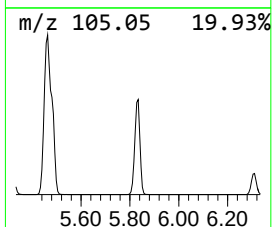
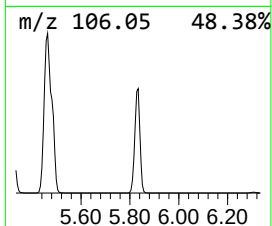
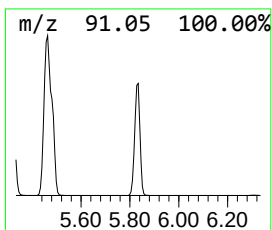
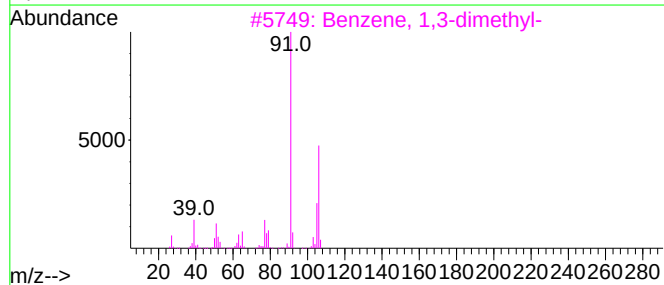
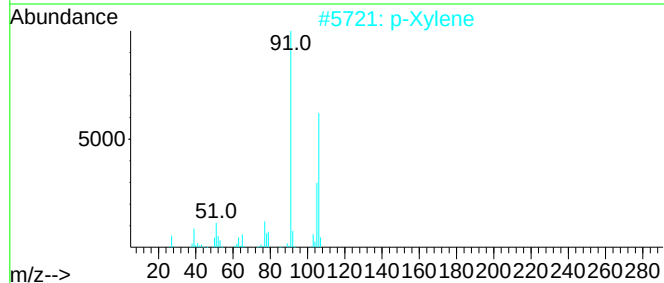
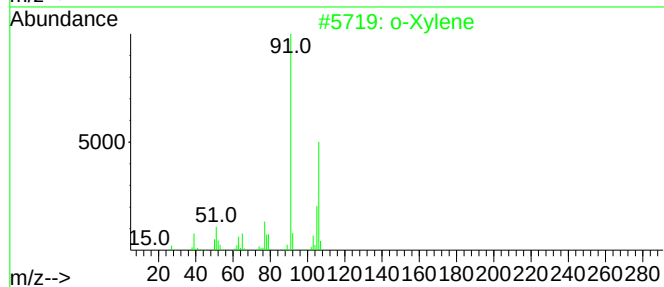
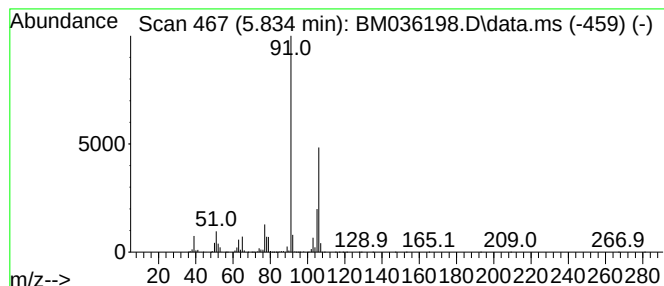
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Peak Number 5 o-Xylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.834	166.10 ng	21932600	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	97
2			p-Xylene	106	C8H10	000106-42-3	97
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4			Ethylbenzene	106	C8H10	000100-41-4	87
5			Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	86



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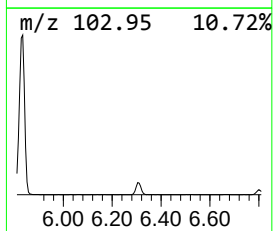
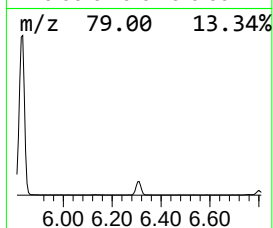
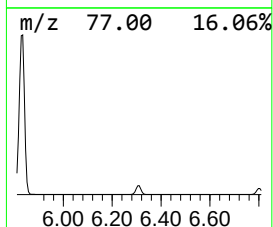
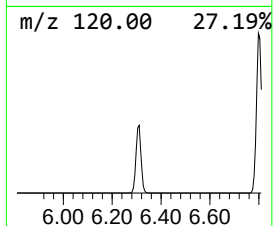
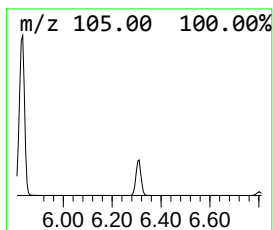
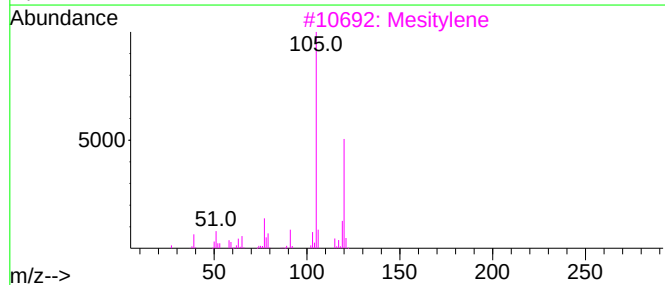
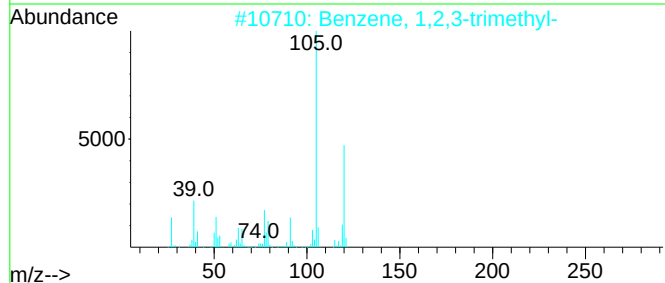
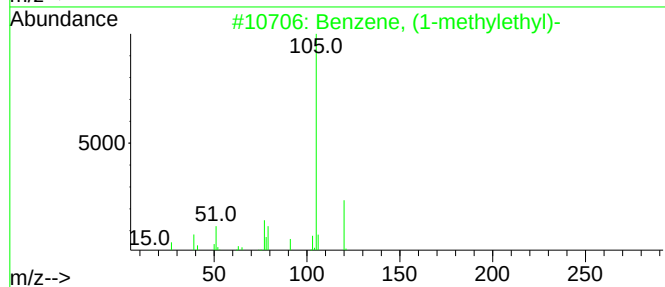
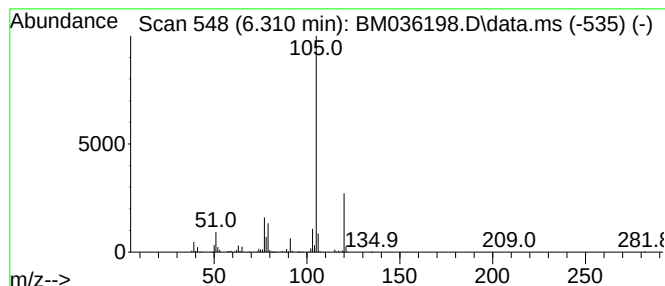
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.310	6.55 ng	864976	1,4-Dichlorobenzene-d4	7.834

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			Mesitylene	120	C9H12	000108-67-8	86
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	83
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72



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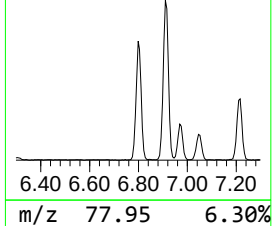
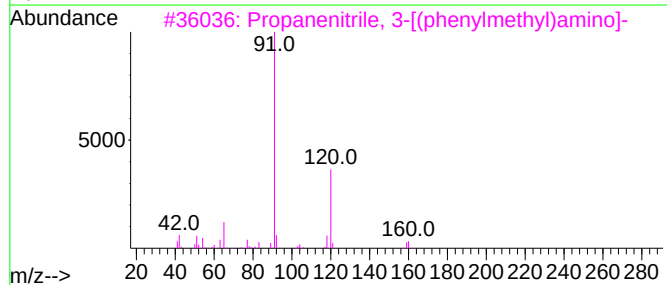
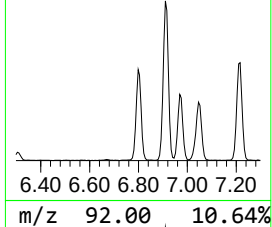
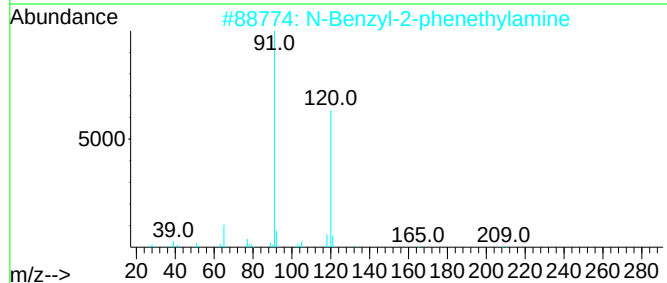
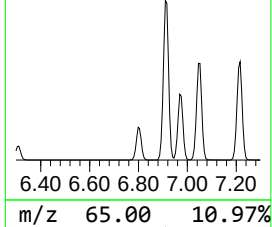
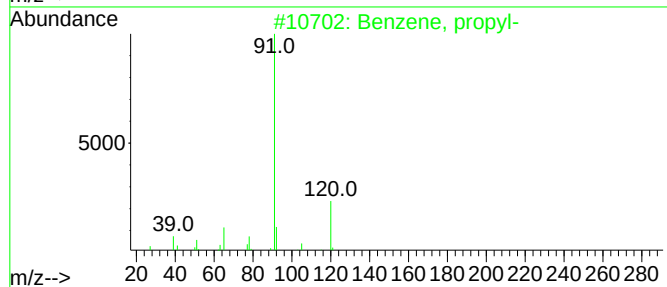
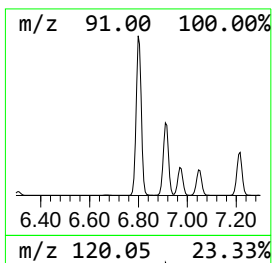
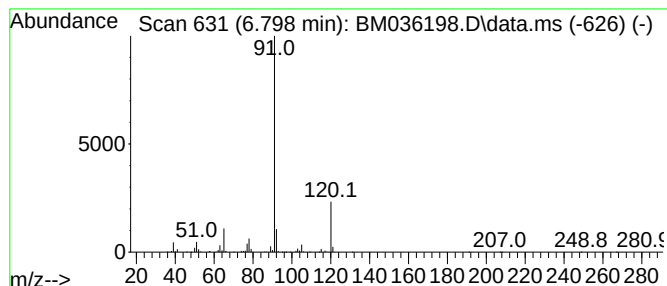
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.798	13.82 ng	1824720	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
3			Propanenitrile, 3-[(phenylmethyl)amino]-	160	C10H12N2	000706-03-6	72
4			1,2-Ethanediamine, N,N'-bis(phenyl)-	240	C16H20N2	000140-28-3	64
5			N-Isobutyl(phenyl)methanesulfonamide	227	C11H17NO2S	144615-94-1	59



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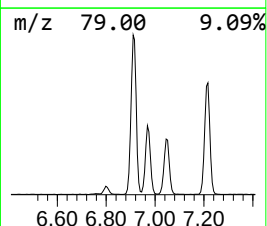
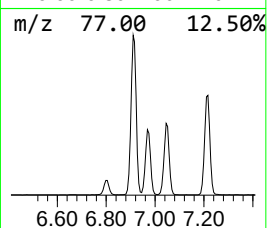
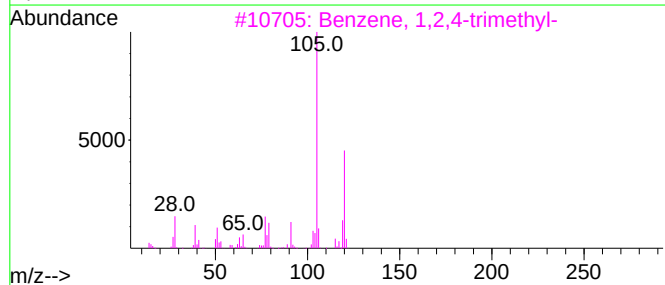
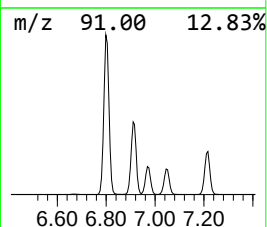
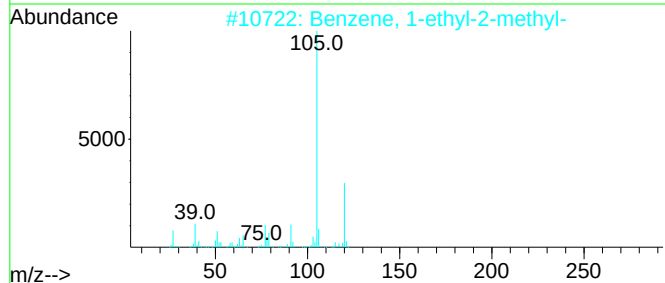
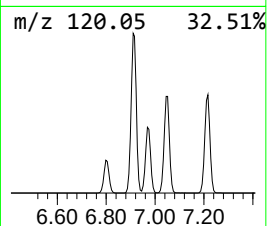
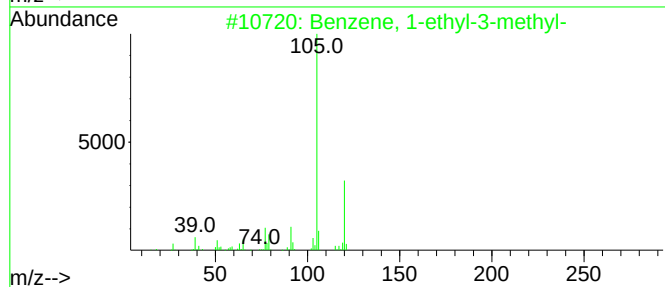
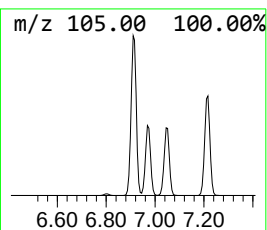
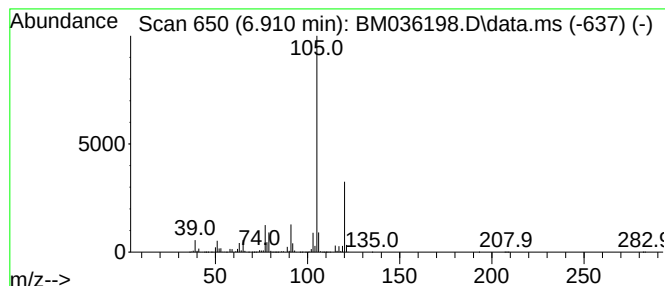
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.910	64.65 ng	8536630	1,4-Dichlorobenzene-d4	7.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
Operator : CG/JU
Sample : N3971-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

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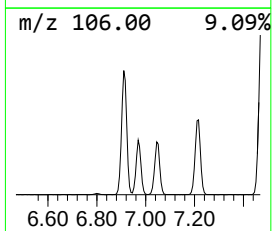
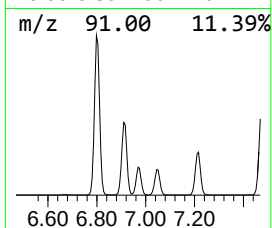
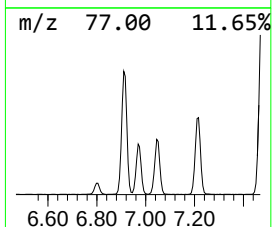
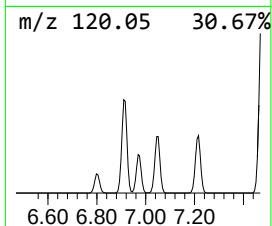
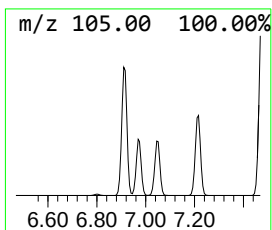
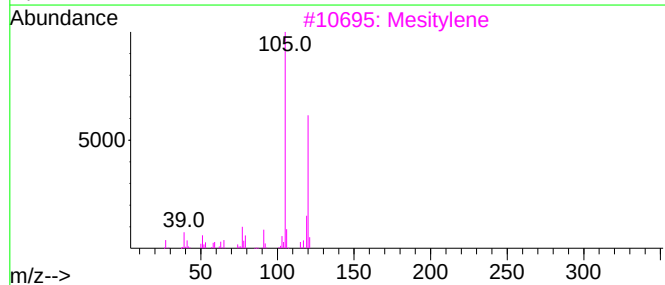
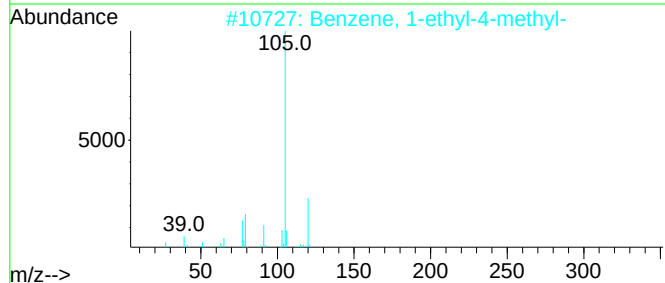
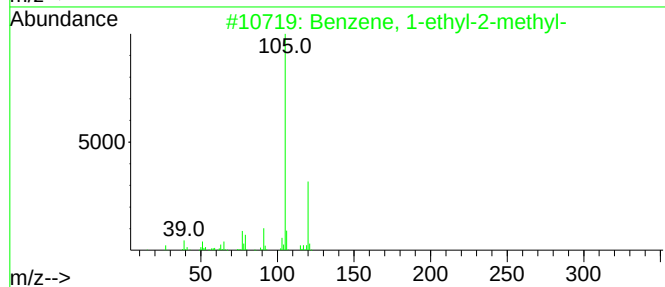
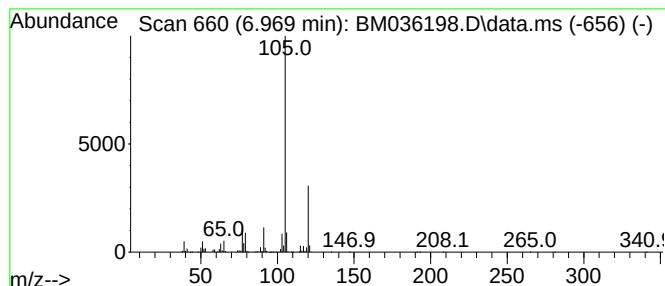
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.969	25.85 ng	3413490	1,4-Dichlorobenzene-d4	7.834

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-4-methyl-	120	C9H12	000622-96-8	94
3			Mesitylene	120	C9H12	000108-67-8	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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
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Misc :
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Instrument :
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ClientSampleId :
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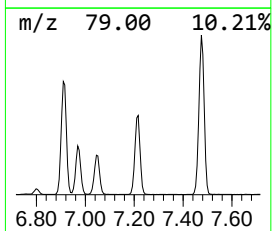
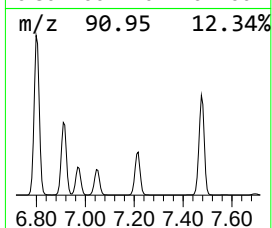
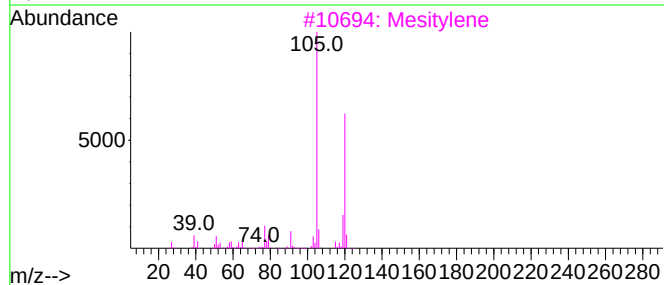
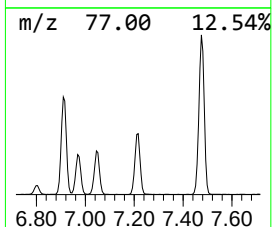
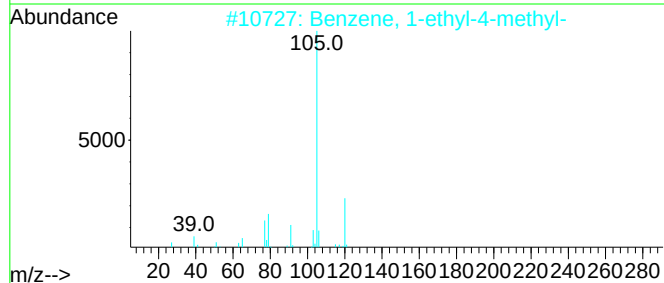
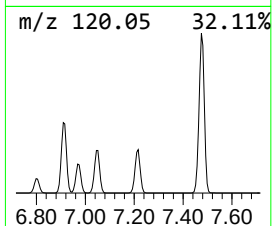
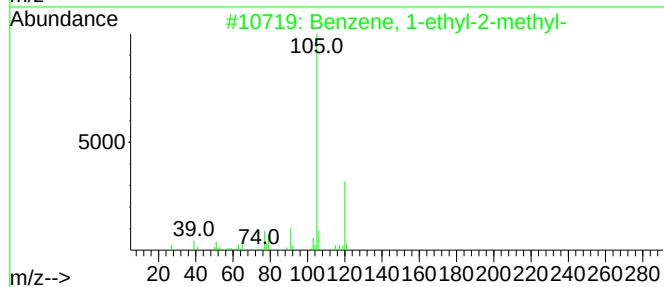
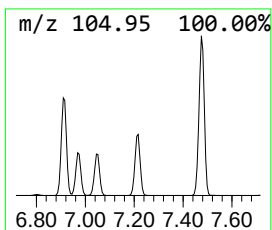
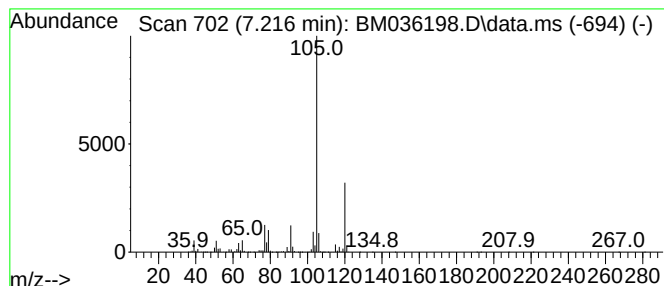
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.216	39.91 ng	5269920	1,4-Dichlorobenzene-d4	7.834

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	93
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
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Misc :
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Instrument :
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ClientSampleId :
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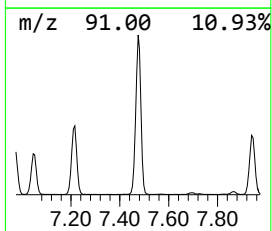
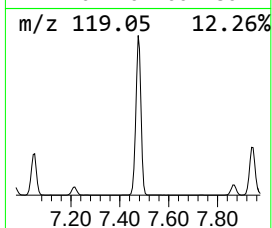
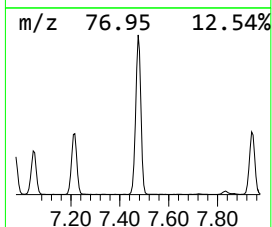
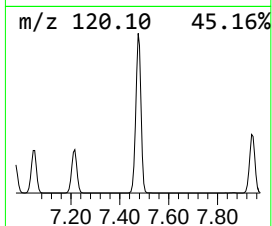
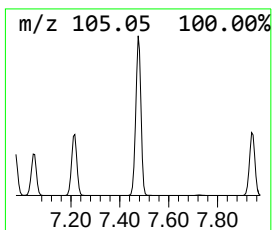
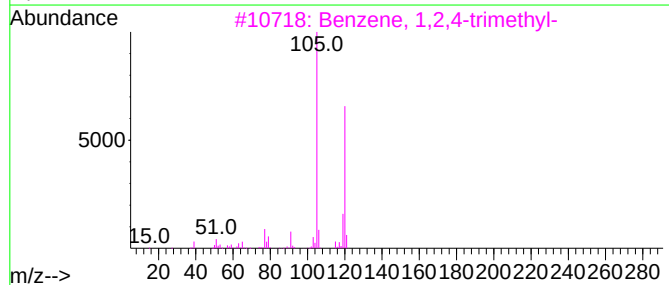
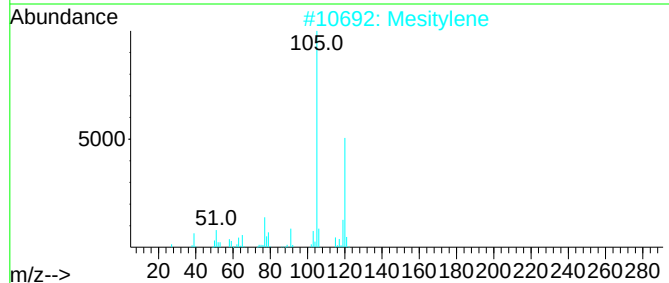
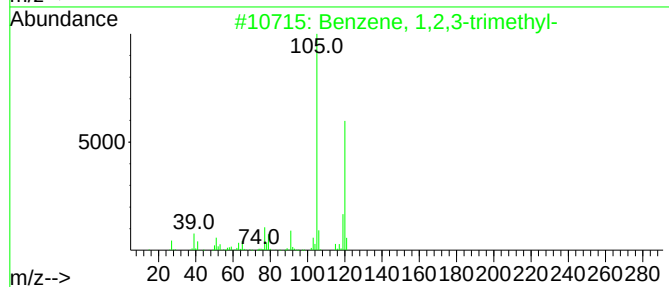
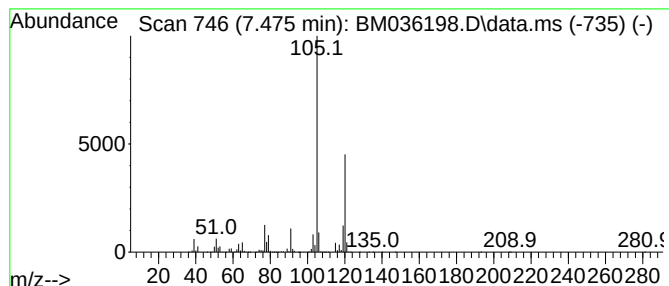
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.475	113.74 ng	15019200	1,4-Dichlorobenzene-d4	7.834

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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
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Misc :
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Instrument :
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ClientSampleId :
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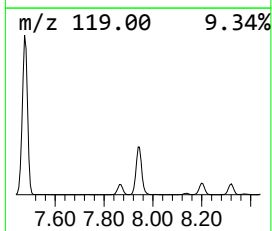
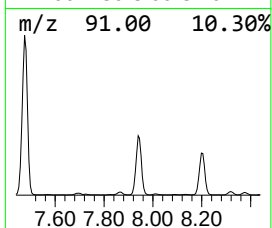
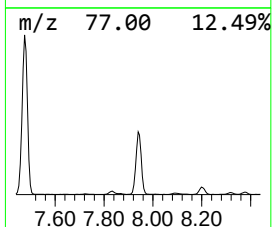
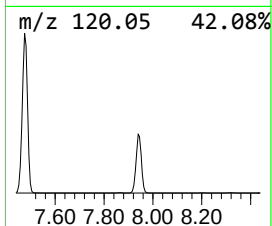
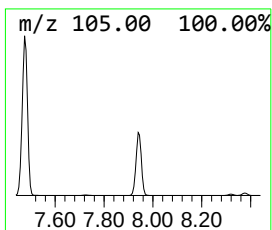
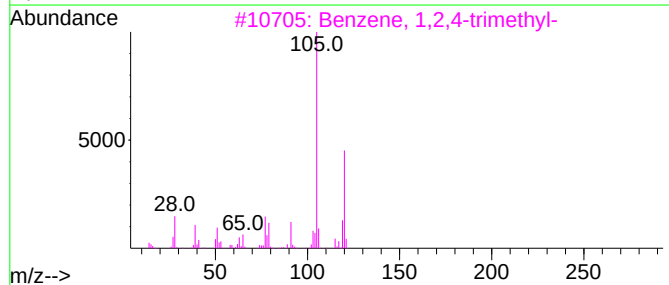
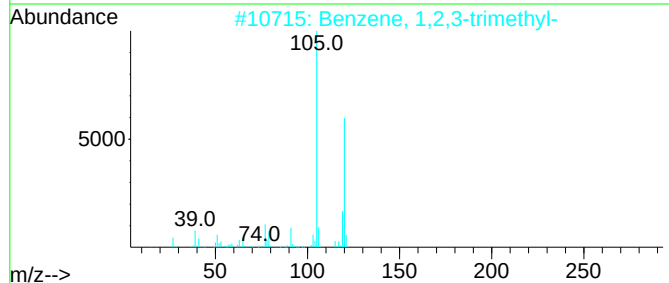
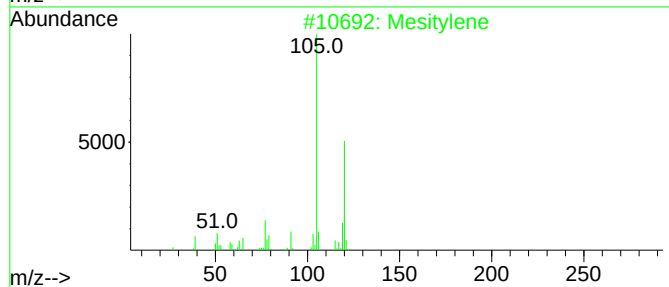
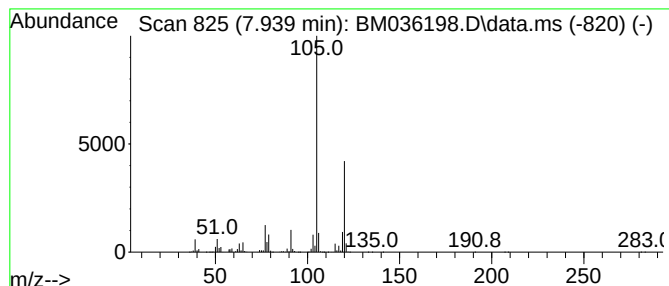
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.939	45.38 ng	5992080	1,4-Dichlorobenzene-d4	7.834

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	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
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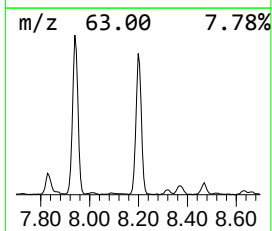
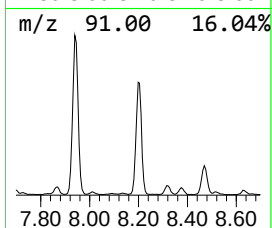
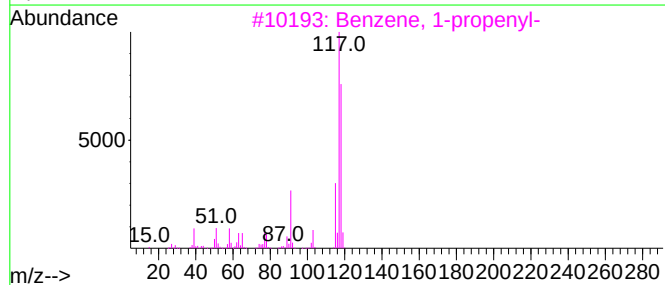
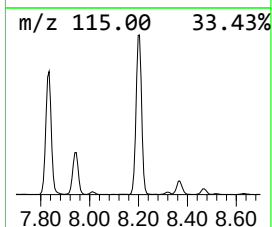
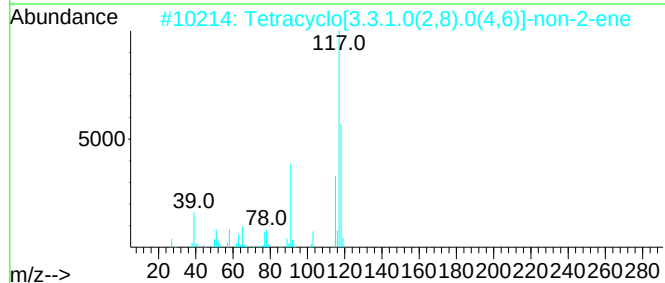
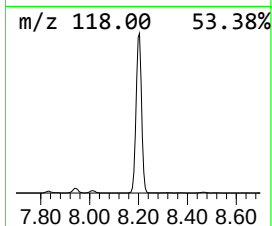
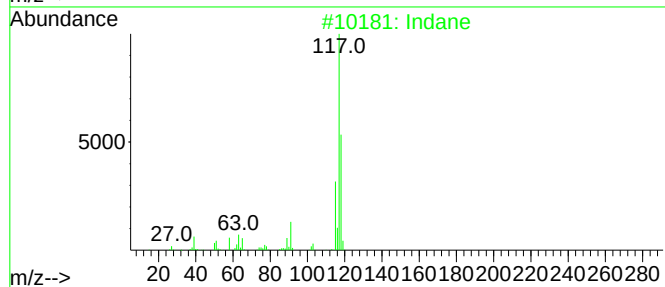
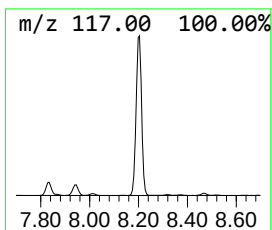
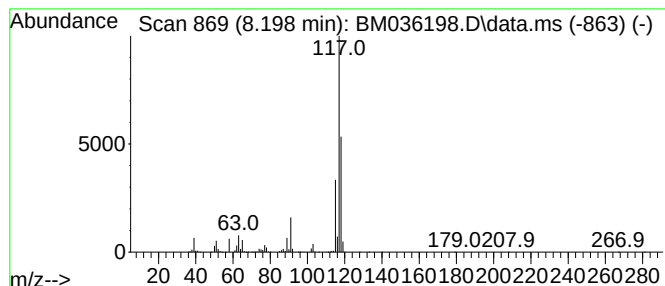
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.198	22.59 ng	2982910	1,4-Dichlorobenzene-d4	7.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	95
2		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
3		Benzene, 1-propenyl-	118	C9H10	000637-50-3	53
4		Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	53
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
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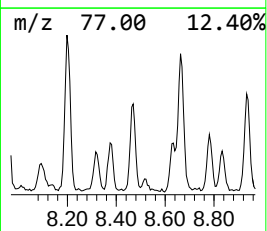
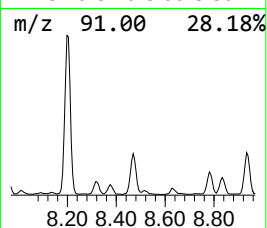
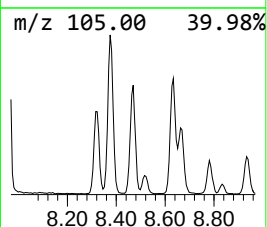
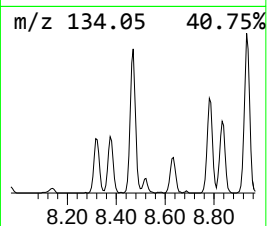
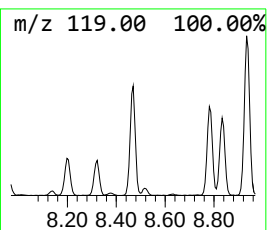
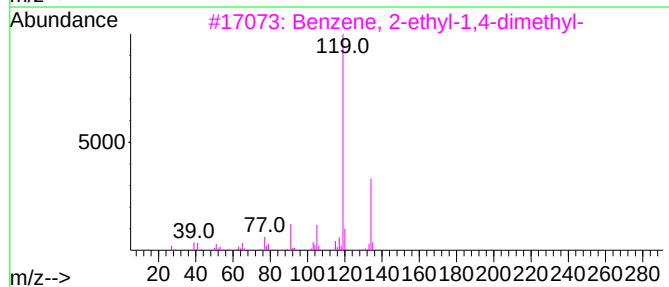
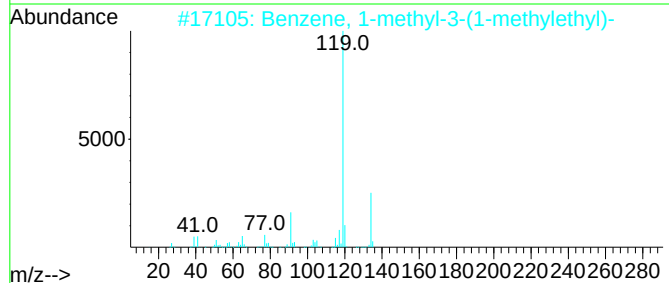
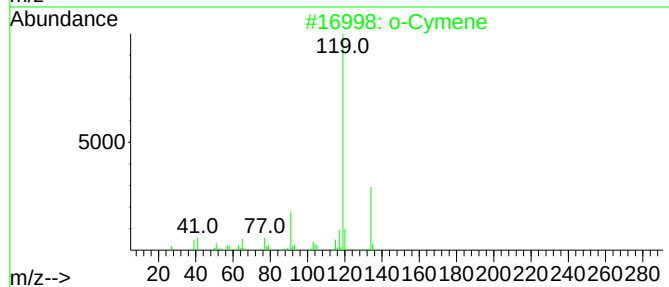
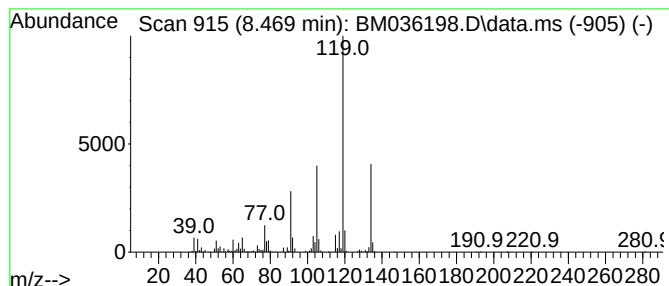
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 o-Cymene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.469	4.93 ng	650817	1,4-Dichlorobenzene-d4	7.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	93
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
Operator : CG/JU
Sample : N3971-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MW-14R

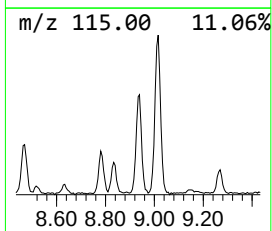
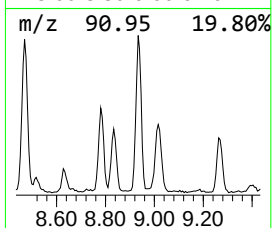
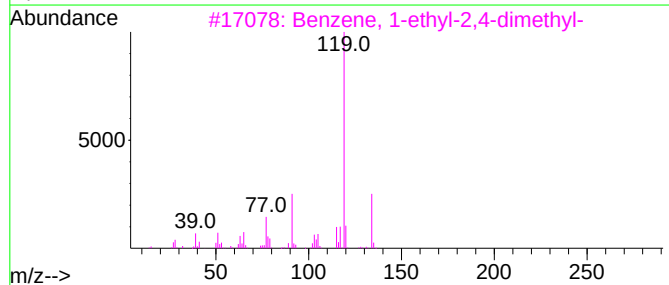
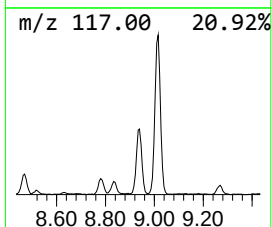
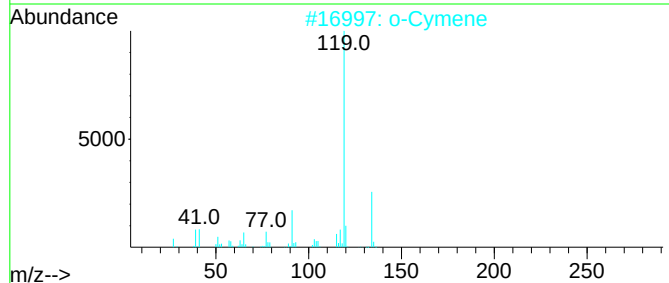
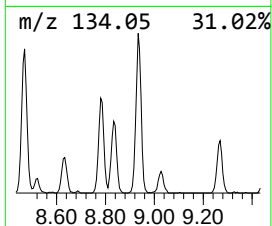
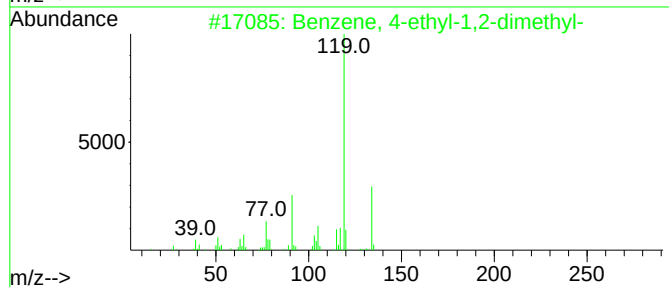
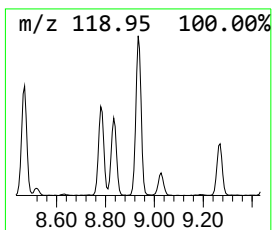
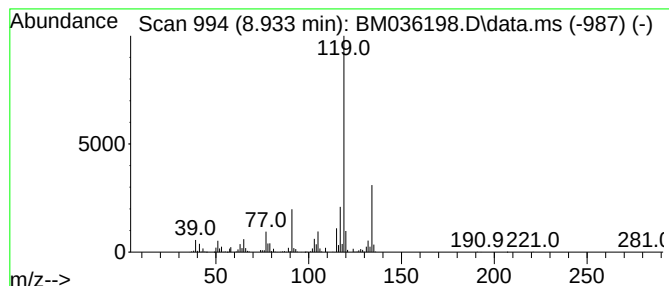
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.933	5.19 ng	684874	1,4-Dichlorobenzene-d4	7.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2		o-Cymene	134	C10H14	000527-84-4	94
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
Operator : CG/JU
Sample : N3971-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

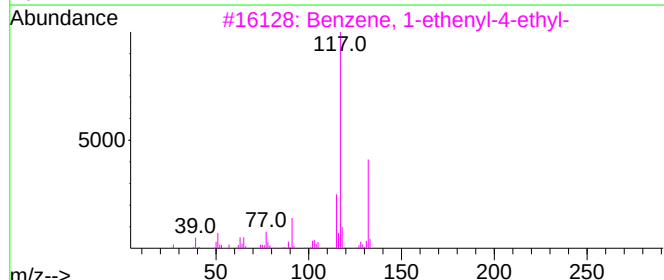
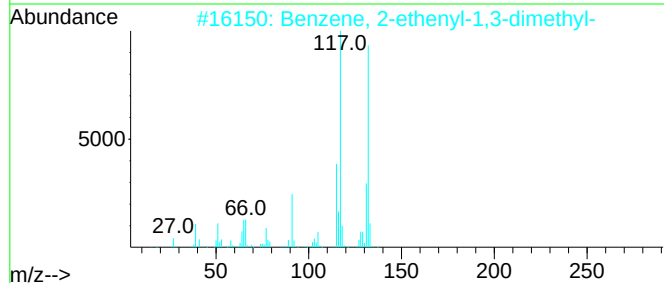
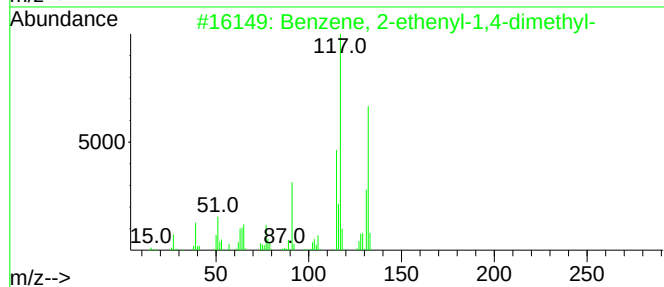
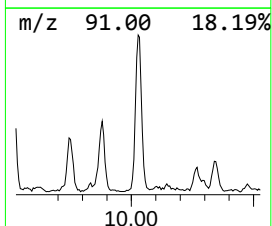
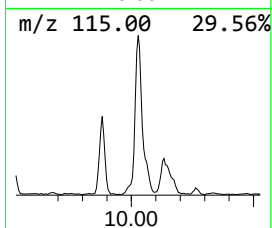
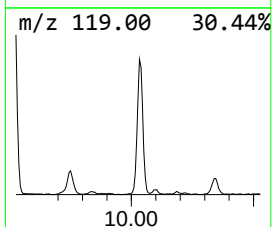
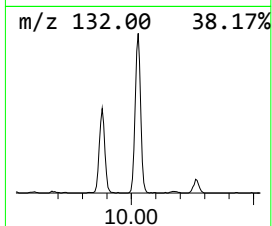
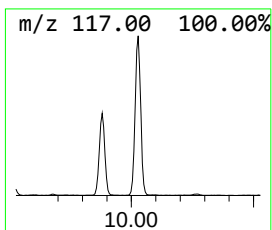
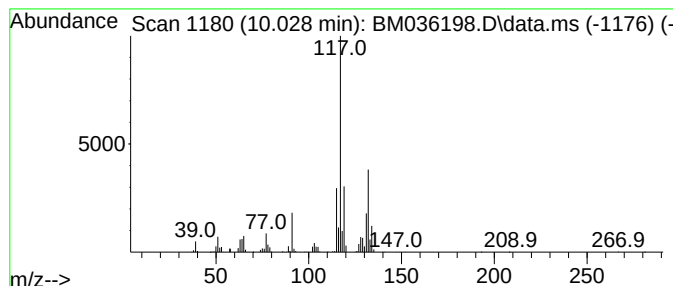
Instrument :
BNA_M
ClientSampleId :
MW-14R

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.028	6.22 ng	1062620	Naphthalene-d8	10.633	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	91
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036198.D
Acq On : 10 Aug 2022 18:46
Operator : CG/JU
Sample : N3971-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

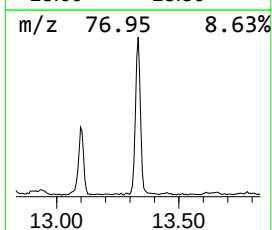
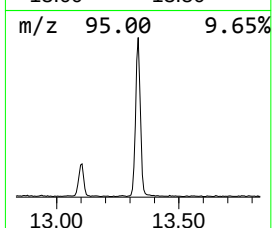
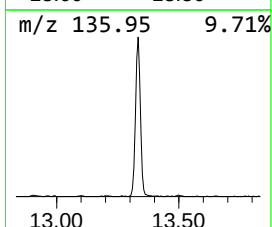
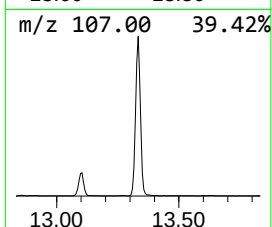
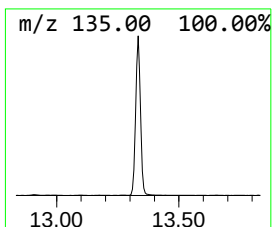
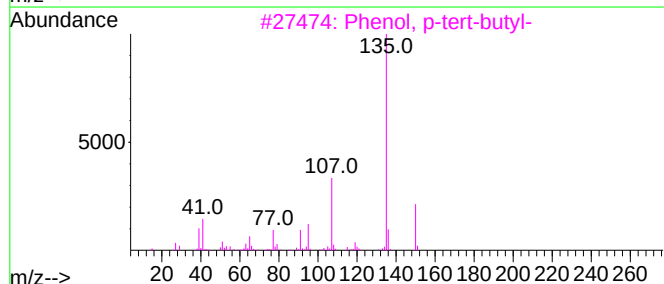
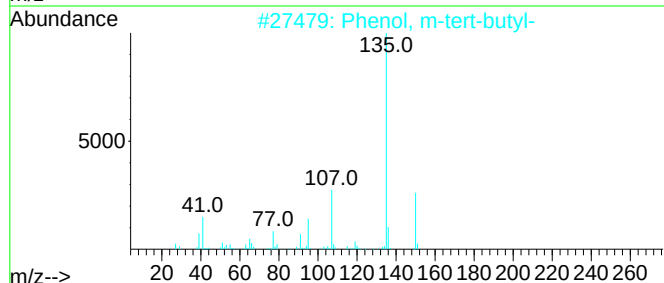
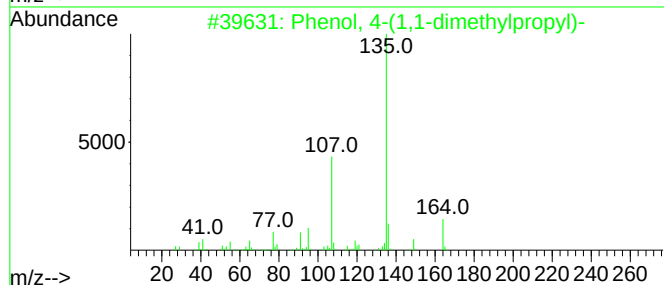
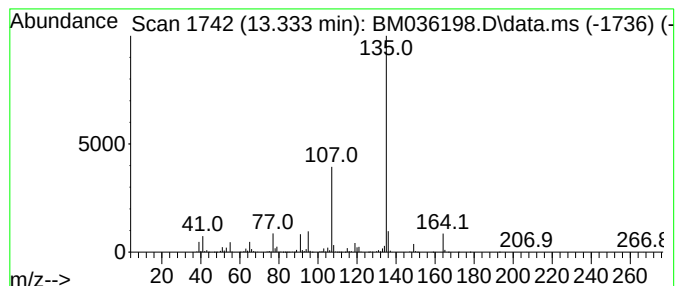
Instrument :
BNA_M
ClientSampleId :
MW-14R

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Phenol, 4-(1,1-dimethylprop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.333	9.89 ng	2089140	Acenaphthene-d10	14.474	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	94
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	86
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
4	Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	64
5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
 Data File : BM036198.D
 Acq On : 10 Aug 2022 18:46
 Operator : CG/JU
 Sample : N3971-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-14R

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.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
Cyclopentanone	4.281	13.1	ng	1724900	1	7.834	2640930	20.0
2-Pentanone, 4-...	4.934	6.0	ng	787242	1	7.834	2640930	20.0
o-Xylene	5.834	166.1	ng	21932600	1	7.834	2640930	20.0
Benzene, (1-met...	6.310	6.5	ng	864976	1	7.834	2640930	20.0
Benzene, propyl-	6.798	13.8	ng	1824720	1	7.834	2640930	20.0
Benzene, 1-ethy...	6.910	64.7	ng	8536630	1	7.834	2640930	20.0
Benzene, 1-ethy...	6.969	25.9	ng	3413490	1	7.834	2640930	20.0
Benzene, 1-ethy...	7.216	39.9	ng	5269920	1	7.834	2640930	20.0
Benzene, 1,2,3-...	7.475	113.7	ng	15019200	1	7.834	2640930	20.0
Mesitylene	7.939	45.4	ng	5992080	1	7.834	2640930	20.0
Indane	8.198	22.6	ng	2982910	1	7.834	2640930	20.0
o-Cymene	8.469	4.9	ng	650817	1	7.834	2640930	20.0
Benzene, 4-ethy...	8.933	5.2	ng	684874	1	7.834	2640930	20.0
Benzene, 2-ethe...	10.028	6.2	ng	1062620	2	10.633	3419160	20.0
Phenol, 4-(1,1-...	13.333	9.9	ng	2089140	3	14.474	4222900	20.0