

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
 Data File : BP009202.D
 Acq On : 17 Feb 2022 18:58
 Operator : CG/JU
 Sample : N1575-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-9R

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\8270E-BP020722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009202.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.975	163	168	172	rBV	504536	729089	3.53%	0.373%
2	4.264	212	217	230	rVB	421680	718082	3.48%	0.367%
3	4.634	275	280	287	rBV	188448	294745	1.43%	0.151%
4	5.287	384	391	400	rBV	6396618	9980112	48.31%	5.100%
5	5.375	400	406	409	rVV	1754471	2833746	13.72%	1.448%
6	5.422	409	414	439	rVB	7619886	19976019	96.70%	10.208%
7	5.787	467	476	493	rVB	2994882	4876980	23.61%	2.492%
8	6.258	549	556	570	rBV	1037711	1710255	8.28%	0.874%
9	6.752	632	640	650	rBV	2950431	4676693	22.64%	2.390%
10	6.857	651	658	663	rVV	1689272	2648458	12.82%	1.353%
11	6.916	663	668	674	rVV	2898073	4534753	21.95%	2.317%
12	6.993	674	681	699	rVV2	3565718	7031791	34.04%	3.593%
13	7.157	703	709	725	rVB	2567538	4156108	20.12%	2.124%
14	7.328	733	738	743	rBV	166564	268898	1.30%	0.137%
15	7.422	743	754	773	rVB	11755441	20657589	100.00%	10.556%
16	7.569	773	779	787	rBV	240267	450531	2.18%	0.230%
17	7.669	793	796	804	rVB	142256	212683	1.03%	0.109%
18	7.775	806	814	818	rBV	751604	1307564	6.33%	0.668%
19	7.887	827	833	842	rVB	4510745	7657092	37.07%	3.913%
20	8.140	870	876	888	rVB	1859459	3026881	14.65%	1.547%
21	8.263	888	897	902	rBV	942986	1497737	7.25%	0.765%
22	8.316	902	906	914	rVB2	184018	299879	1.45%	0.153%
23	8.410	914	922	927	rBV	1412363	2257643	10.93%	1.154%
24	8.457	927	930	942	rVB	202051	327846	1.59%	0.168%
25	8.575	943	950	956	rBV	434429	695218	3.37%	0.355%
26	8.722	968	975	980	rBV	987831	1547251	7.49%	0.791%
27	8.775	980	984	991	rVB	732194	1107961	5.36%	0.566%
28	8.875	994	1001	1006	rBV	2375129	3769140	18.25%	1.926%
29	8.940	1006	1012	1036	rVB2	2927495	6007762	29.08%	3.070%
30	9.204	1051	1057	1065	rVB	554793	944939	4.57%	0.483%
31	9.399	1085	1090	1095	rBV	1044073	1596849	7.73%	0.816%
32	9.457	1095	1100	1109	rVB	1230033	1985245	9.61%	1.014%
33	9.816	1155	1161	1176	rVB	850048	1476128	7.15%	0.754%
34	9.969	1176	1187	1196	rBV3	1763620	3481498	16.85%	1.779%
35	10.569	1278	1289	1293	rBV	1067641	2010381	9.73%	1.027%
36	10.616	1293	1297	1305	rVV	2240135	4106270	19.88%	2.098%

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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\8270E-BP020722.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.234	1563	1572	1581	rBV	904988	1590195	7.70%	0.813%
38	12.457	1598	1610	1630	rVB	726842	1447816	7.01%	0.740%
39	12.775	1654	1664	1679	rBV	318668	619674	3.00%	0.317%
40	13.040	1701	1709	1721	rBV	8727933	13911857	67.35%	7.109%
41	13.269	1739	1748	1765	rVB	1190991	2062687	9.99%	1.054%
42	14.416	1937	1943	1952	rBV	1755968	2796395	13.54%	1.429%
43	15.245	2076	2084	2097	rVB2	216561	380816	1.84%	0.195%
44	15.792	2172	2177	2187	rVB2	141138	229311	1.11%	0.117%
45	15.922	2188	2199	2216	rBV	6901703	11010896	53.30%	5.627%
46	17.180	2406	2413	2428	rBV	2175404	3478351	16.84%	1.777%
47	17.845	2521	2526	2533	rBV2	274822	356414	1.73%	0.182%
48	19.745	2843	2849	2867	rBV	15191668	20087051	97.24%	10.265%
49	20.986	3056	3060	3068	rVB3	149895	252789	1.22%	0.129%
50	21.280	3105	3110	3122	rBV	2420177	3576861	17.31%	1.828%
51	23.668	3510	3516	3535	rVB2	1274868	3029180	14.66%	1.548%

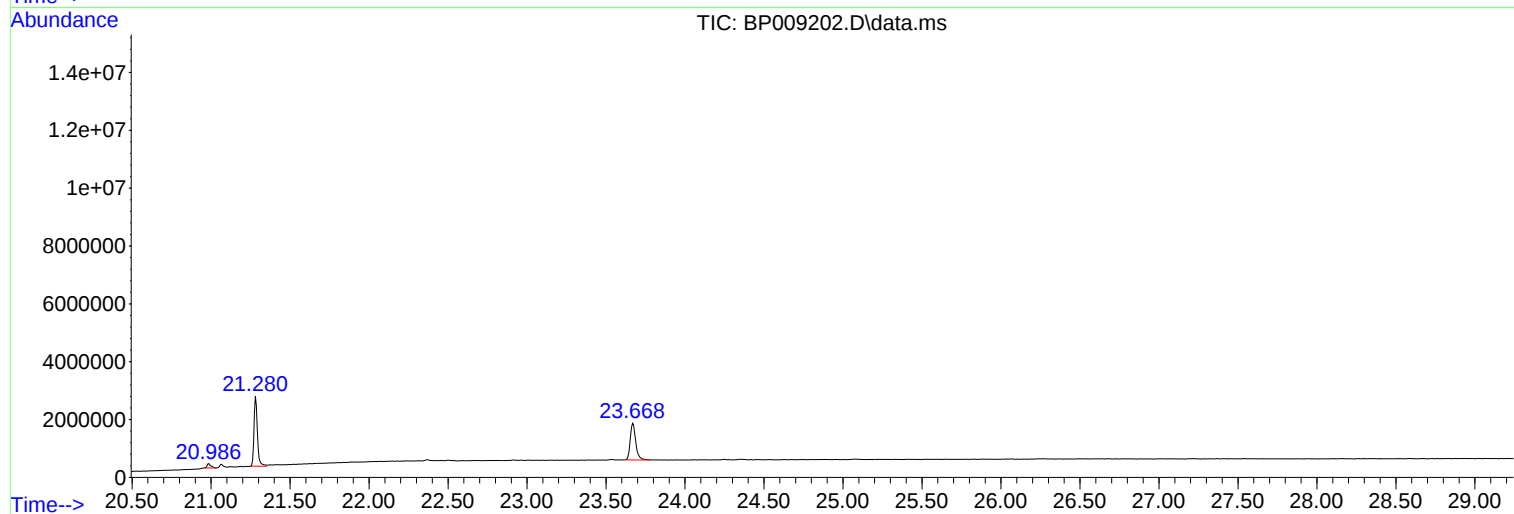
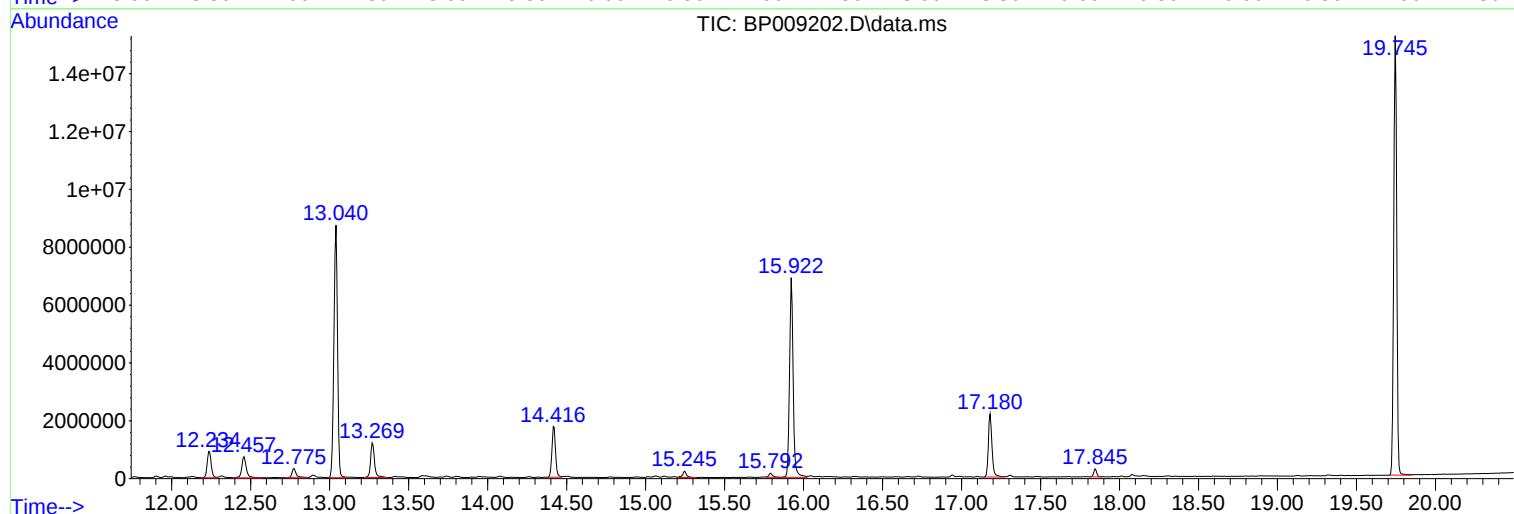
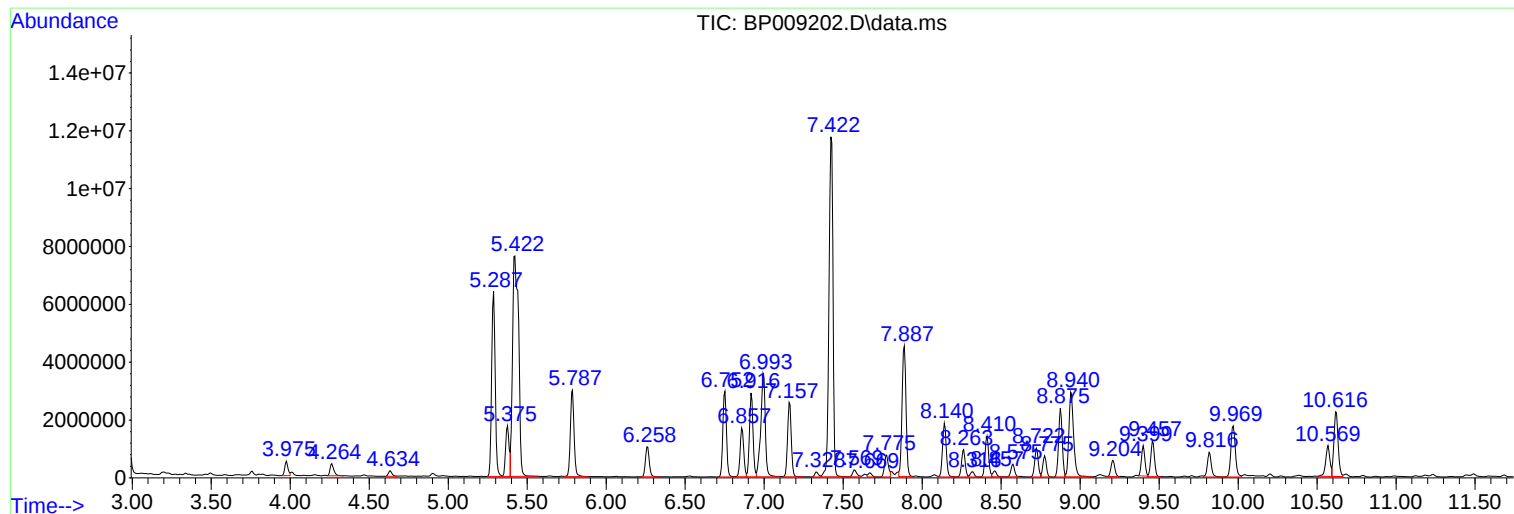
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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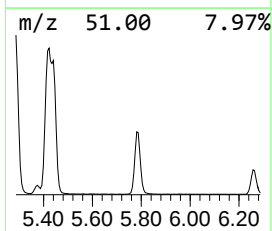
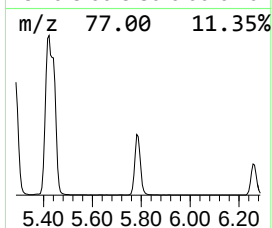
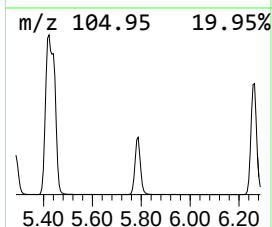
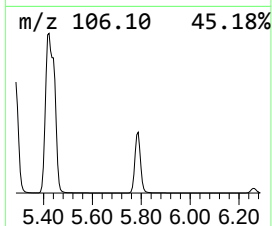
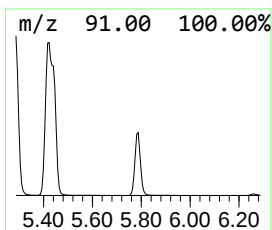
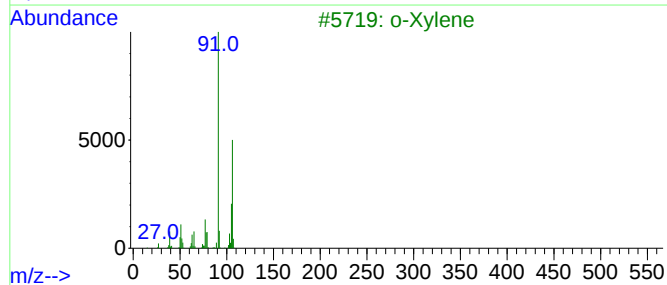
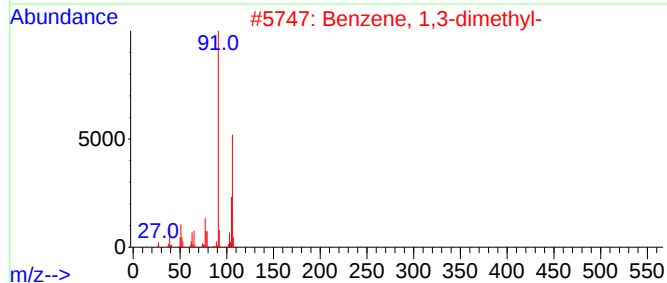
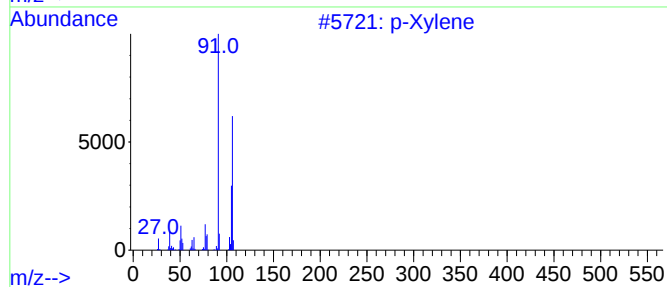
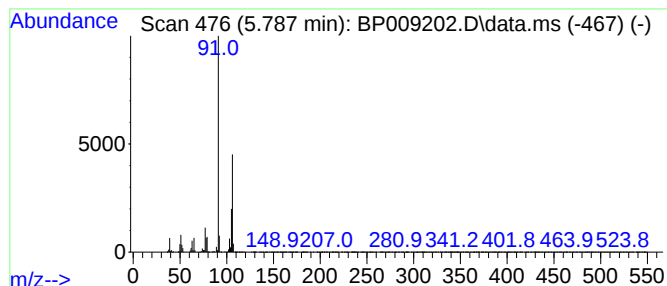
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 p-Xylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.787	74.60 ng	4876980	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Xylene	106	C8H10	000106-42-3	97
2			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3			o-Xylene	106	C8H10	000095-47-6	97
4			Ethylbenzene	106	C8H10	000100-41-4	87
5			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	87



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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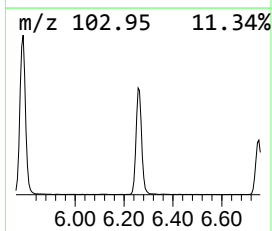
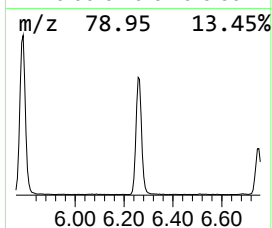
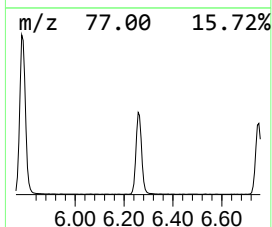
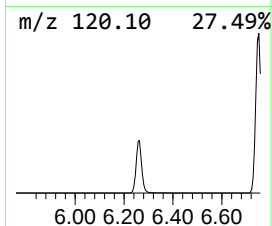
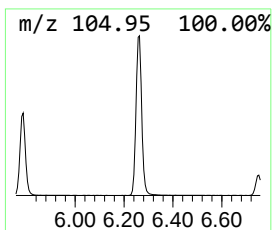
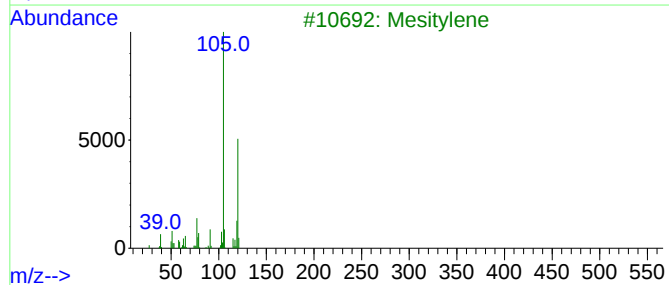
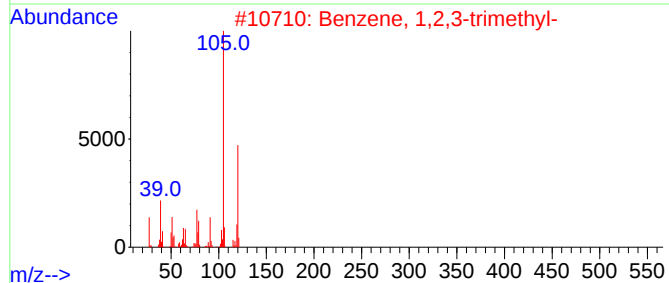
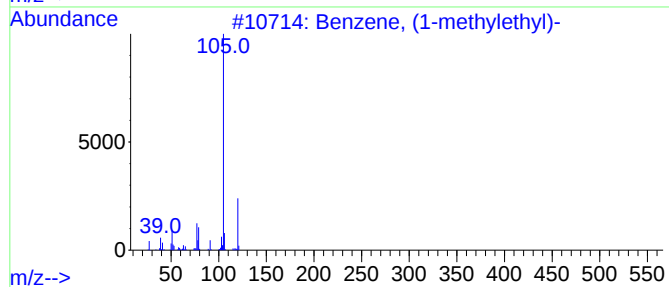
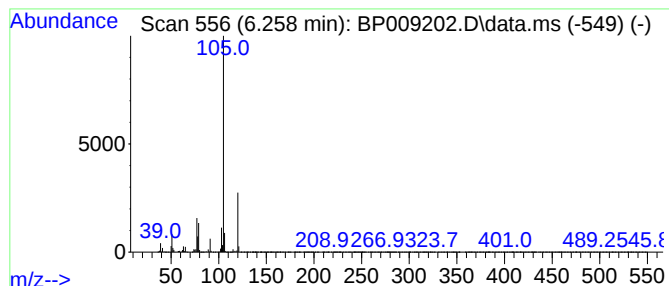
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.258	26.16 ng	1710260	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80



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ALS Vial : 12 Sample Multiplier: 1

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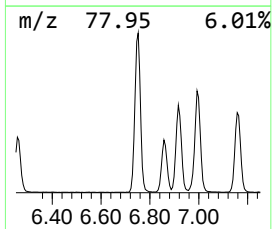
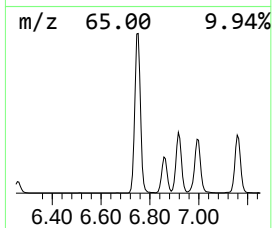
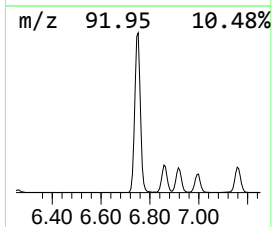
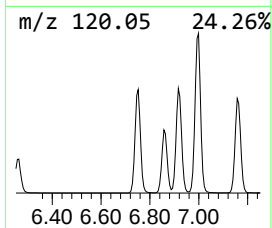
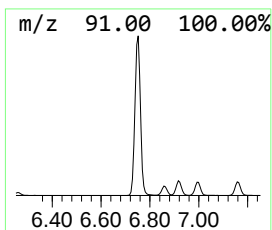
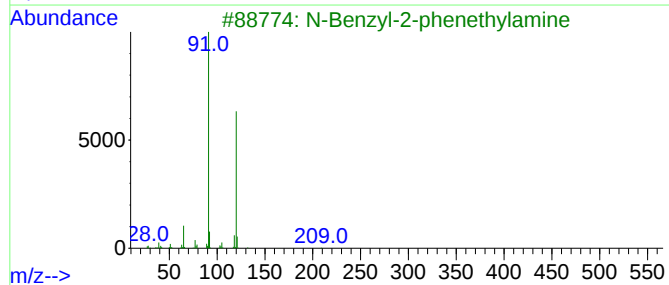
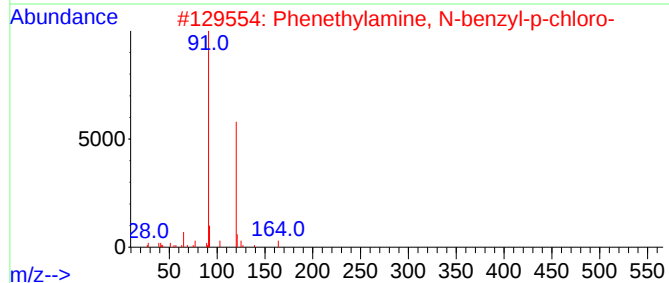
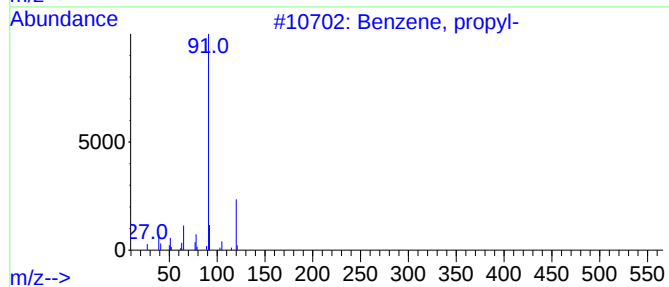
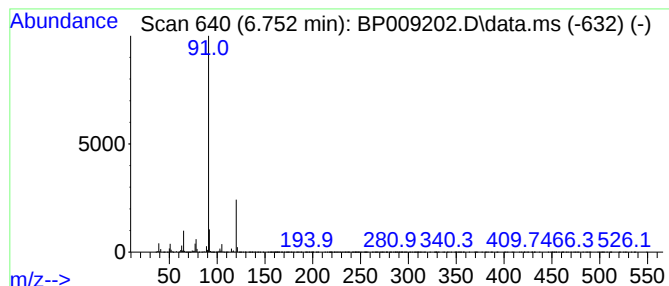
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.752	71.53 ng	4676690	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	91
2			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
3			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
4			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
5			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64



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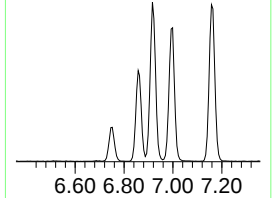
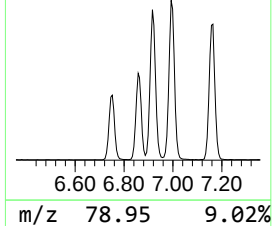
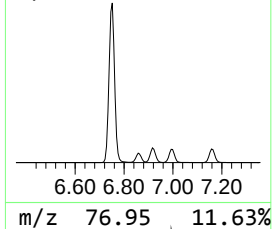
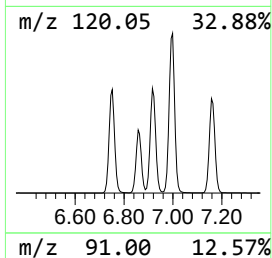
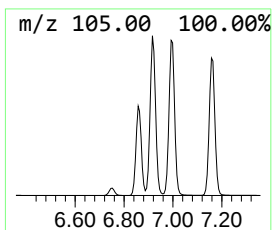
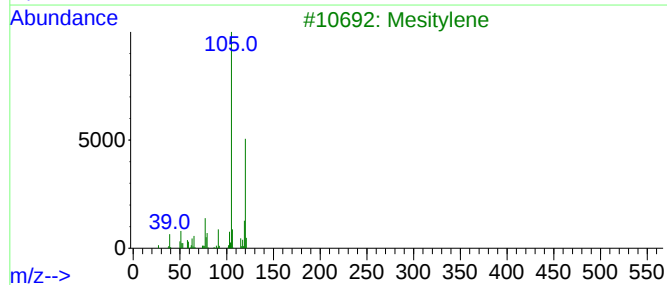
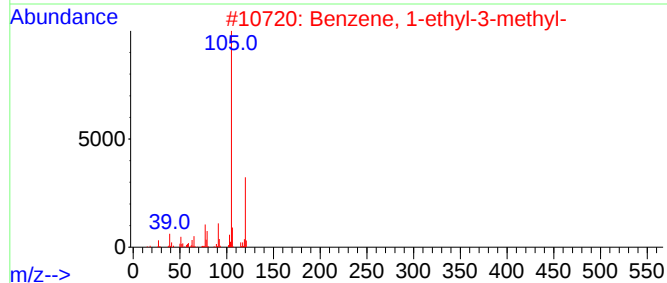
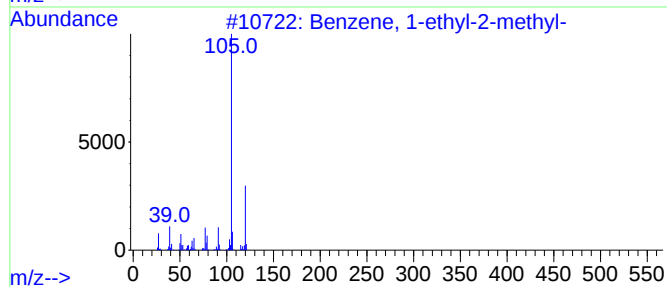
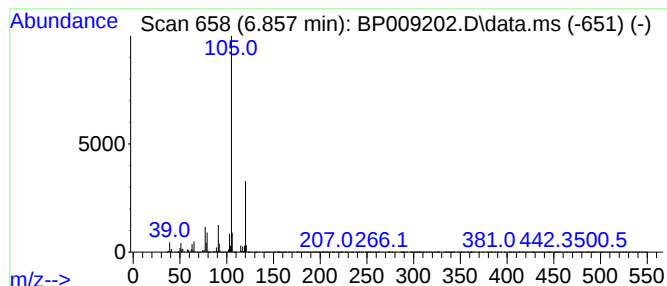
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.857	40.51 ng	2648460	1,4-Dichlorobenzene-d4	7.775

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	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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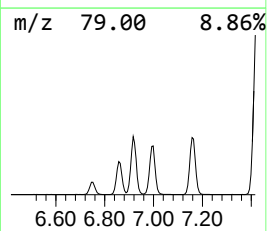
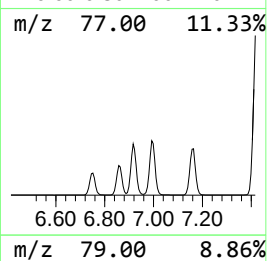
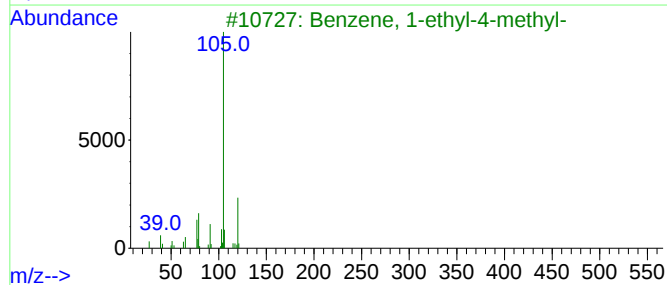
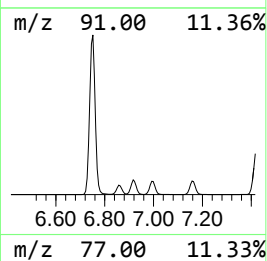
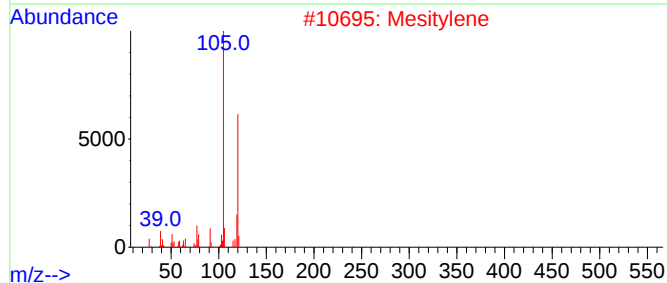
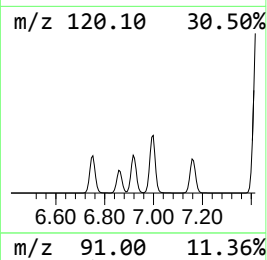
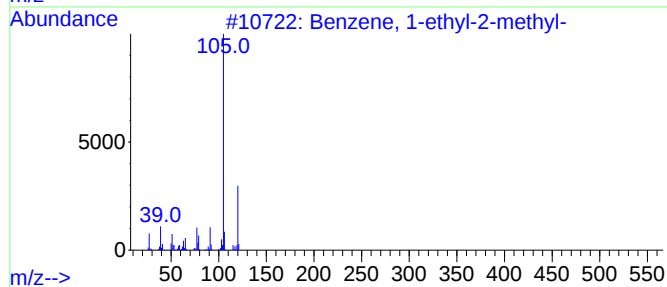
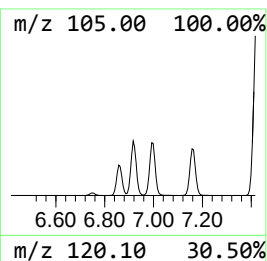
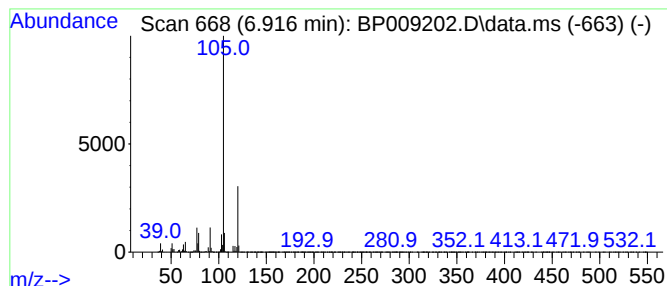
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.916	69.36 ng	4534750	1,4-Dichlorobenzene-d4	7.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-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_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

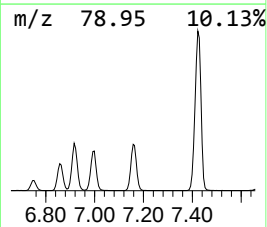
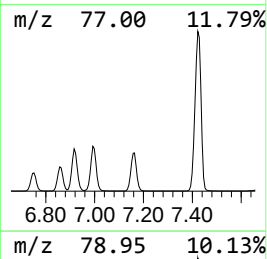
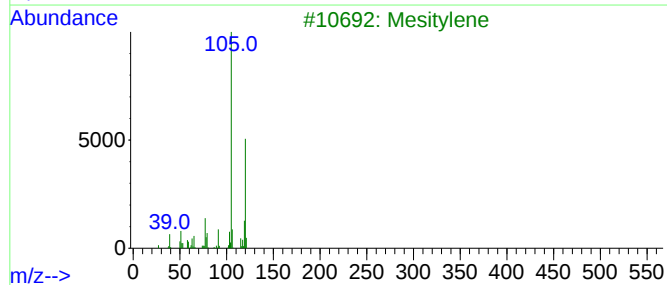
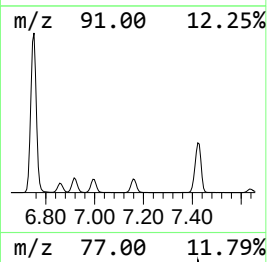
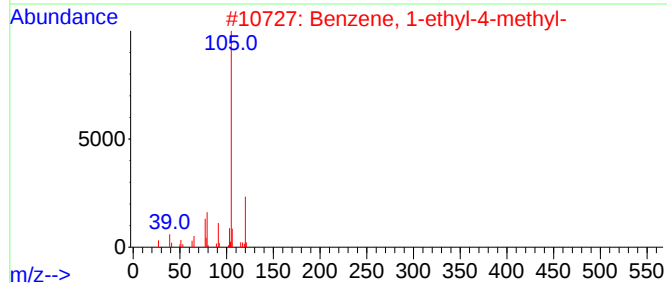
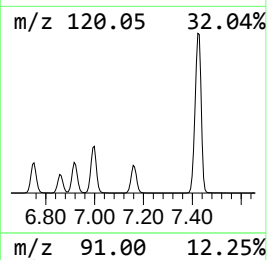
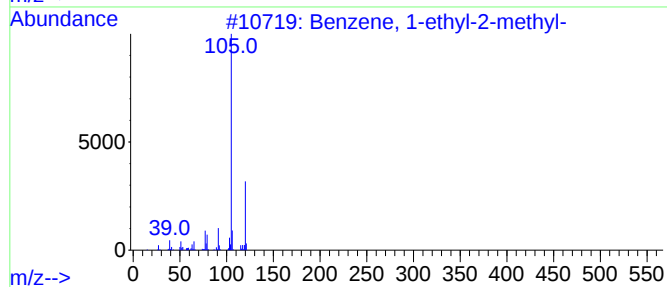
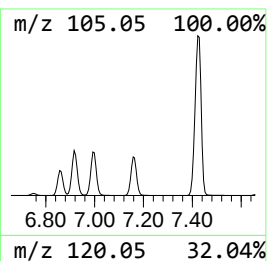
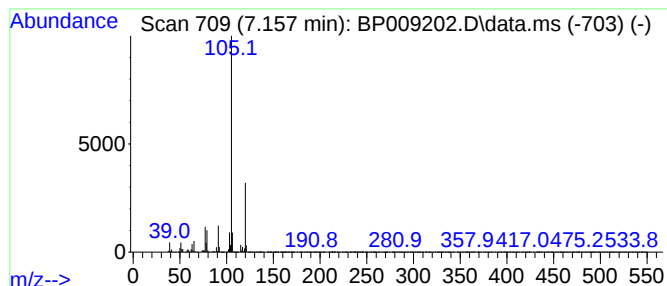
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.157	63.57 ng	4156110	1,4-Dichlorobenzene-d4	7.775

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			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

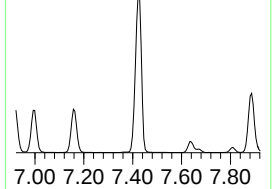
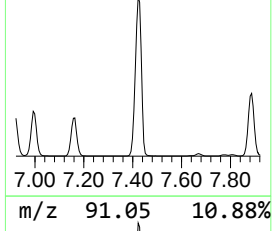
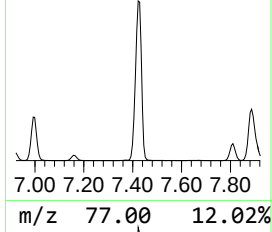
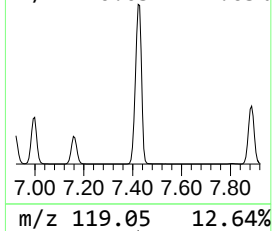
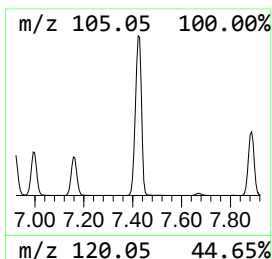
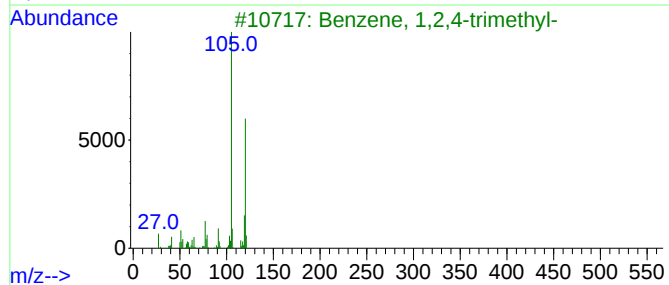
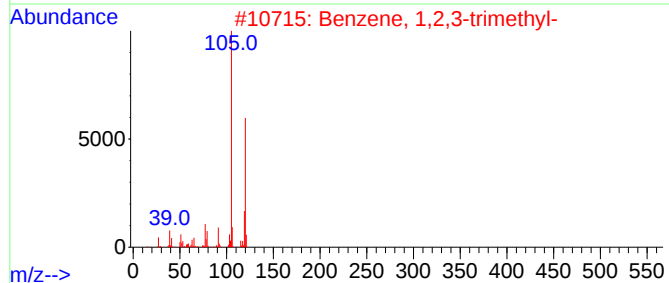
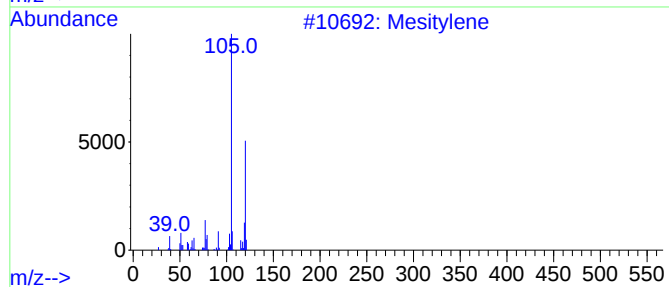
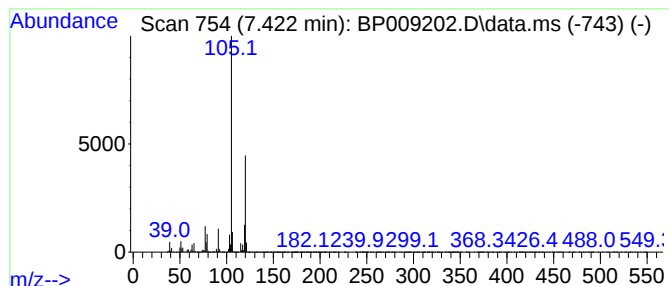
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.422	315.97 ng	20657600	1,4-Dichlorobenzene-d4	7.775

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	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

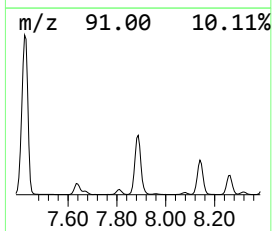
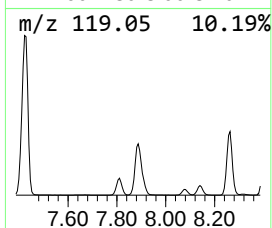
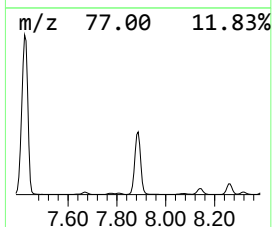
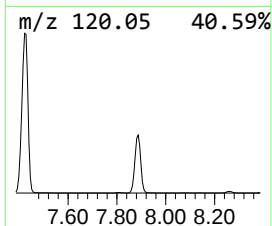
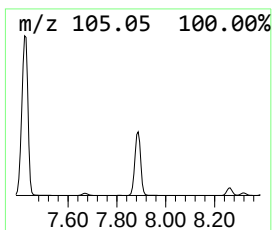
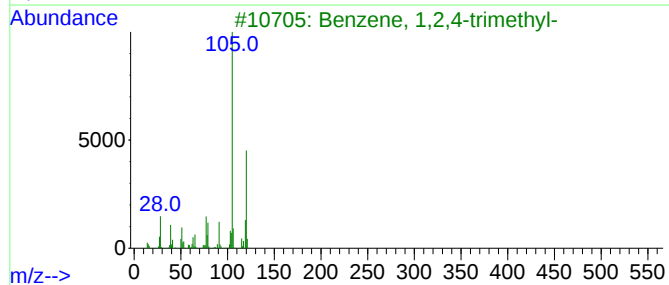
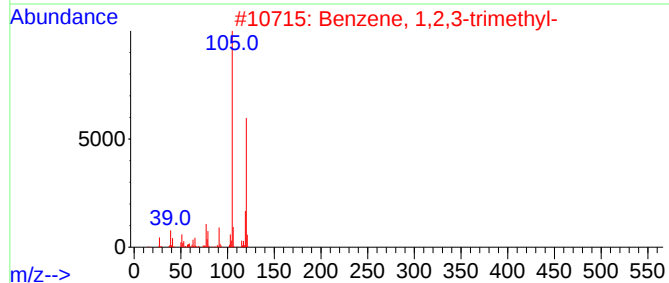
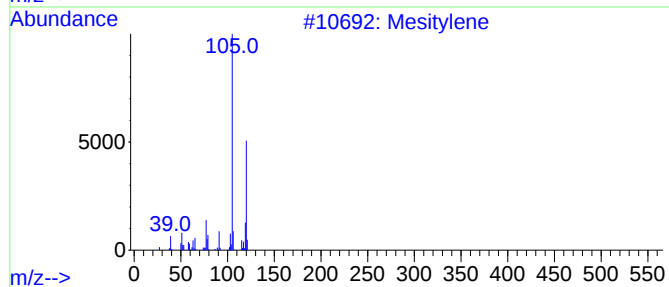
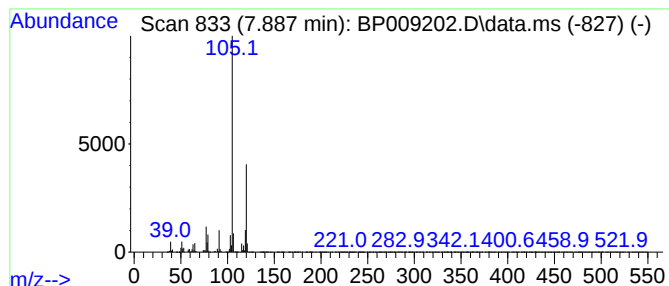
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	117.12 ng	7657090	1,4-Dichlorobenzene-d4	7.775

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-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_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

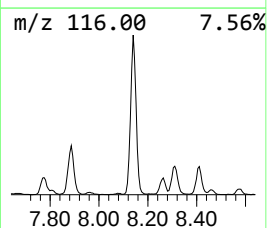
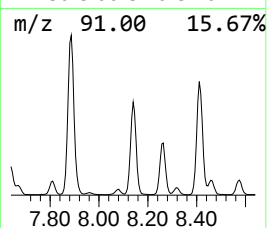
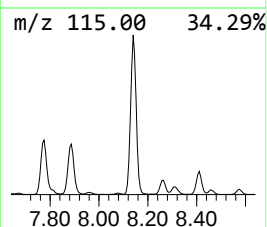
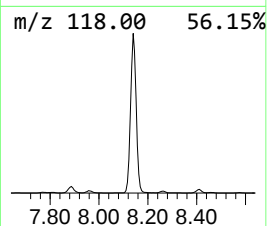
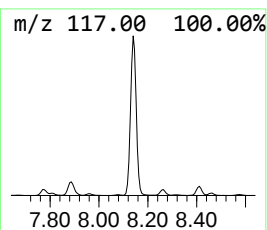
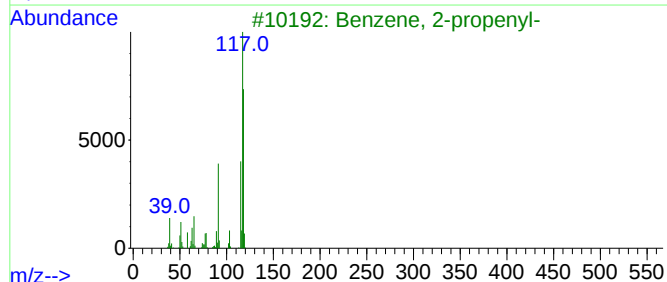
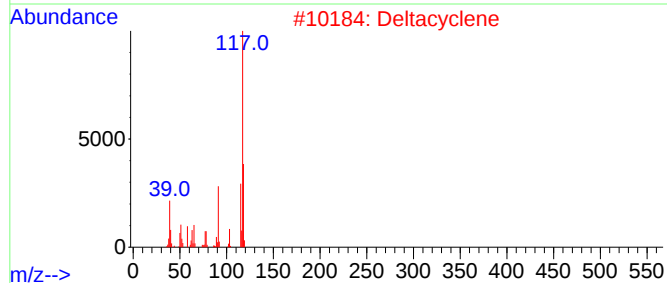
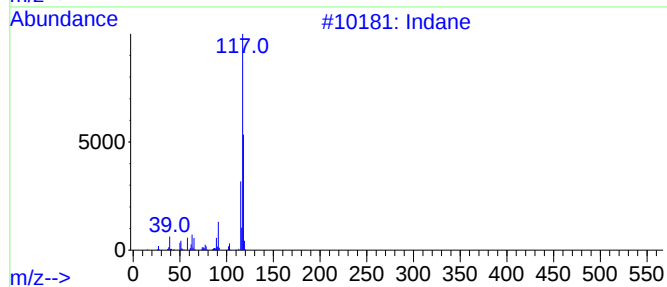
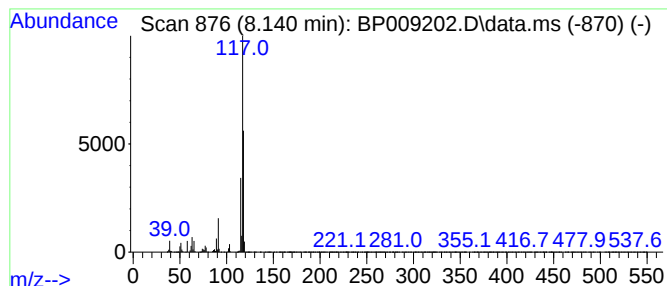
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.140	46.30 ng	3026880	1,4-Dichlorobenzene-d4	7.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	93
2		Deltacyclene	118	C9H10	007785-10-6	72
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

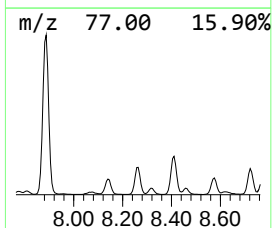
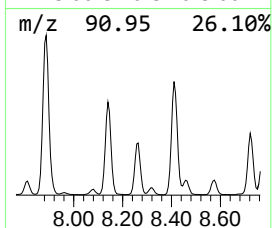
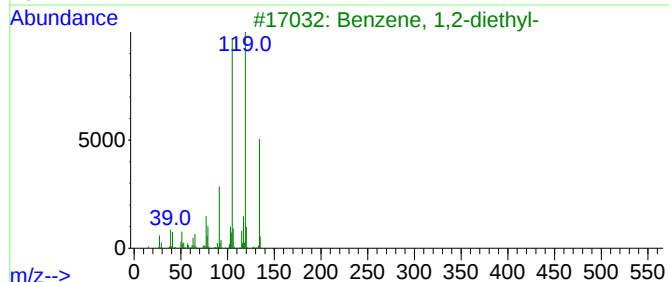
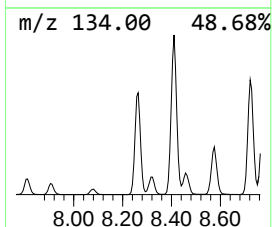
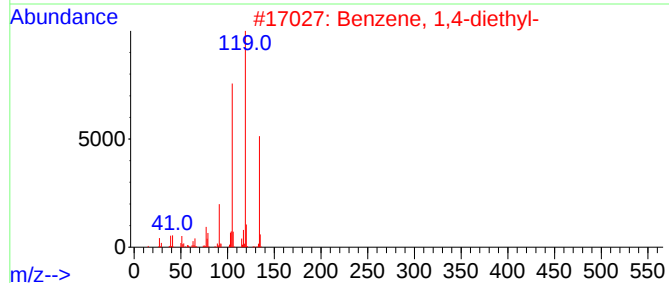
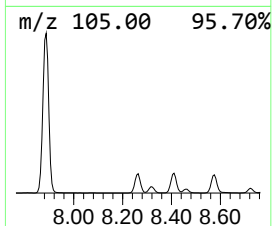
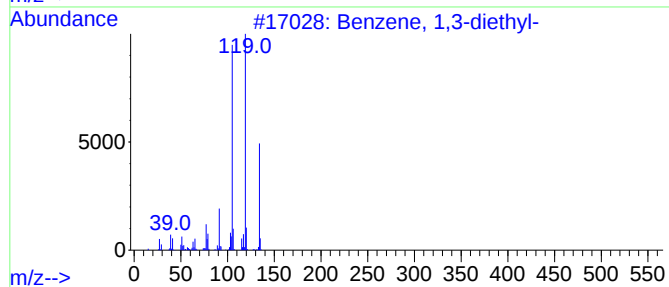
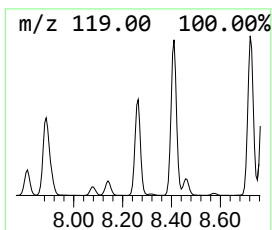
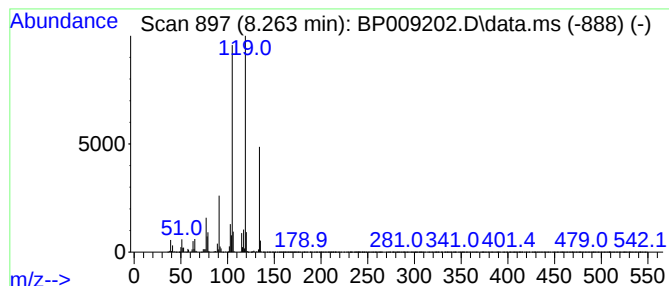
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,3-diethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.263	22.91 ng	1497740	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

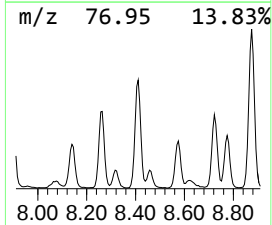
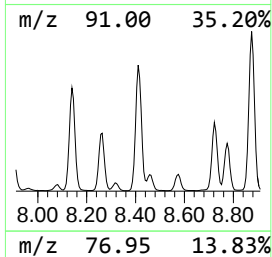
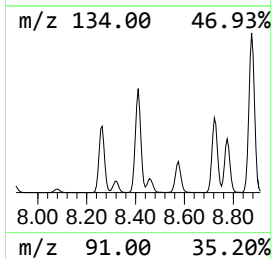
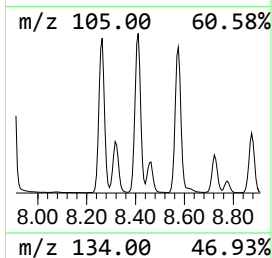
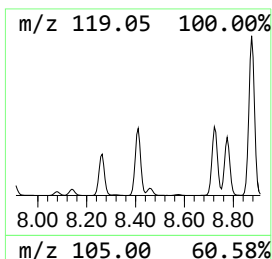
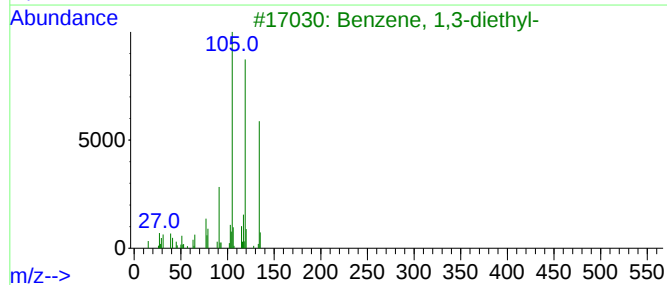
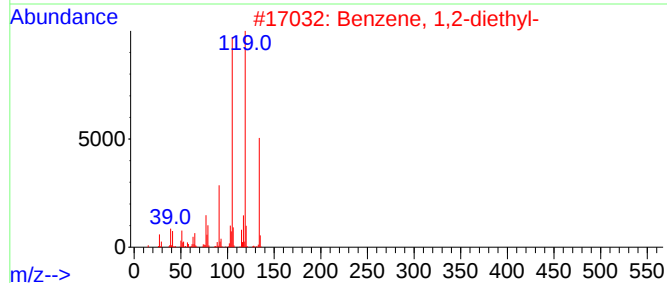
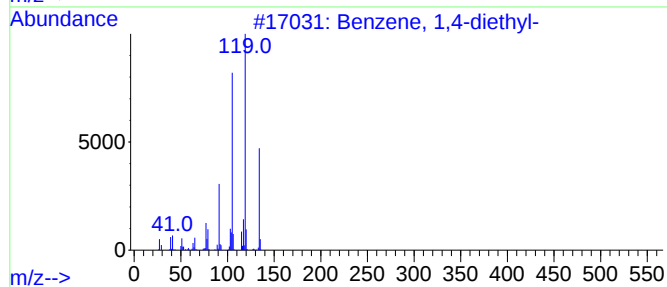
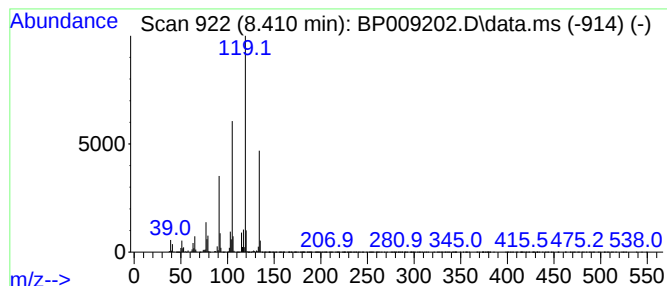
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1,4-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.410	34.53 ng	2257640	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

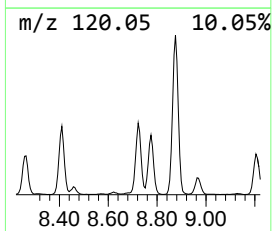
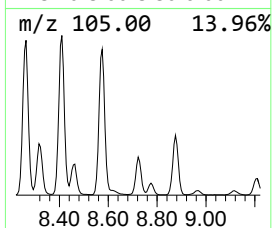
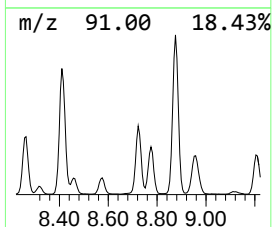
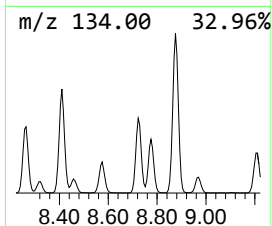
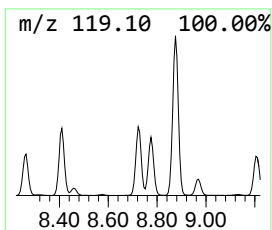
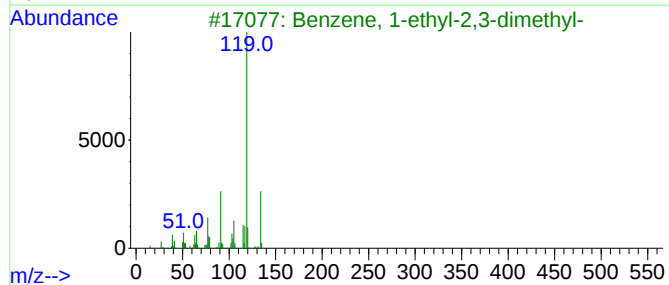
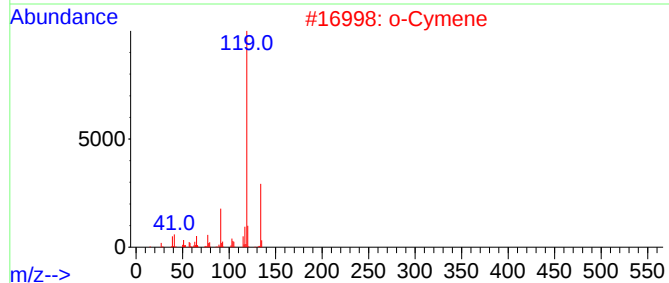
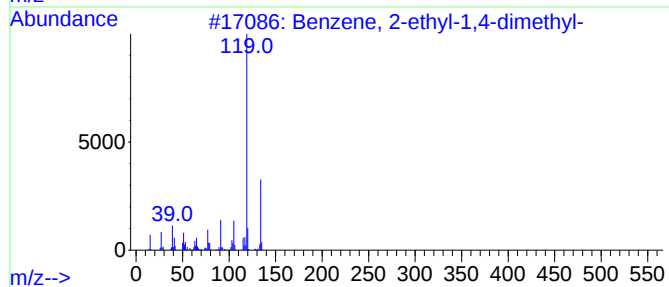
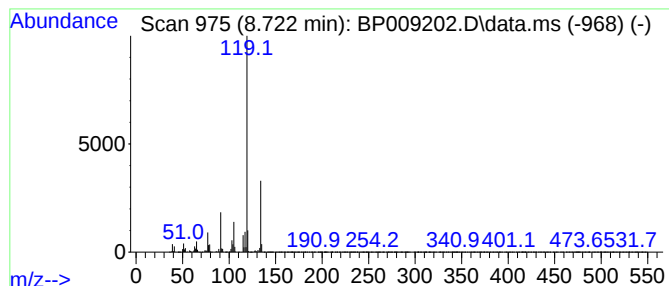
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	23.67 ng	1547250	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-9R

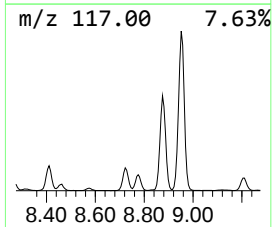
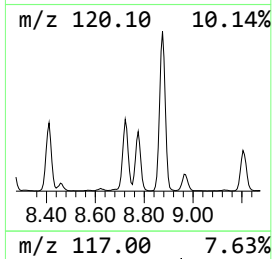
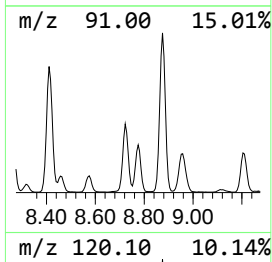
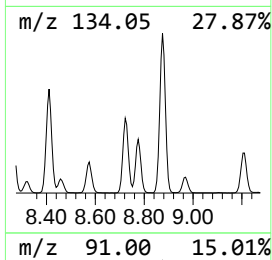
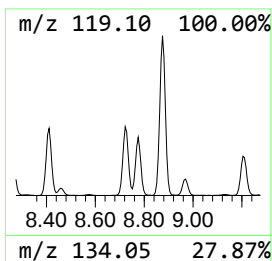
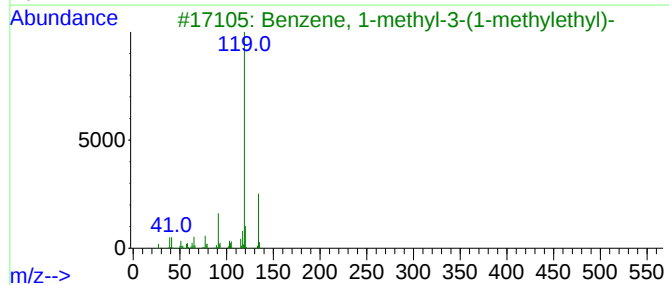
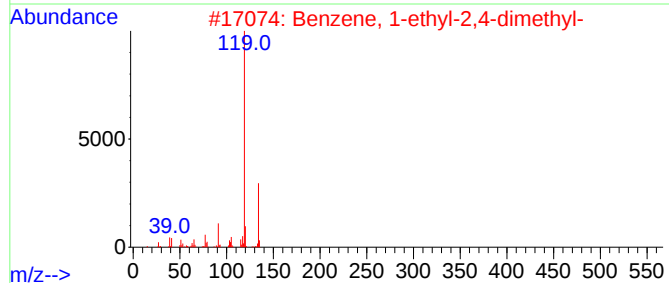
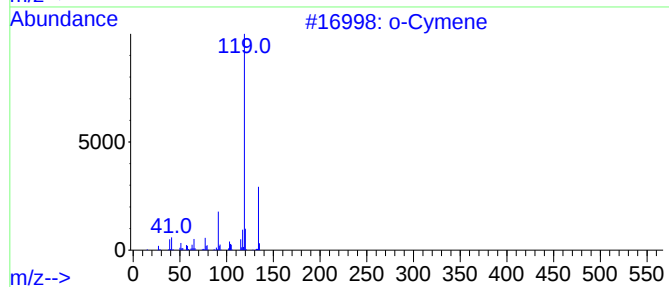
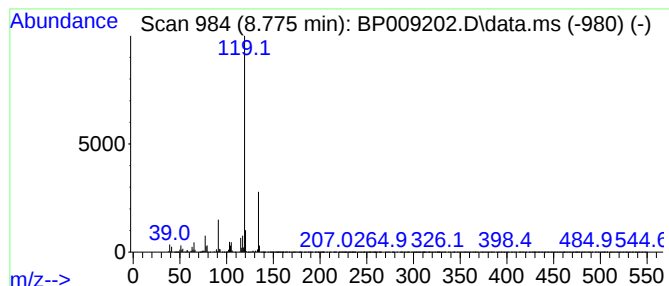
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
8.775	16.95 ng	1107960	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
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_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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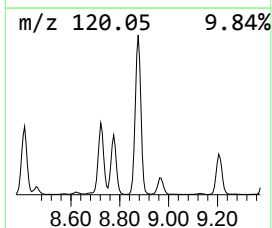
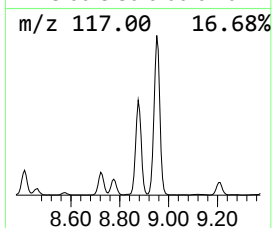
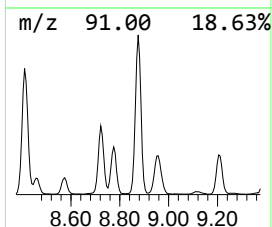
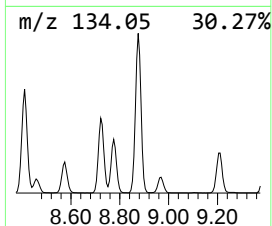
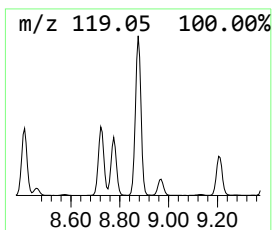
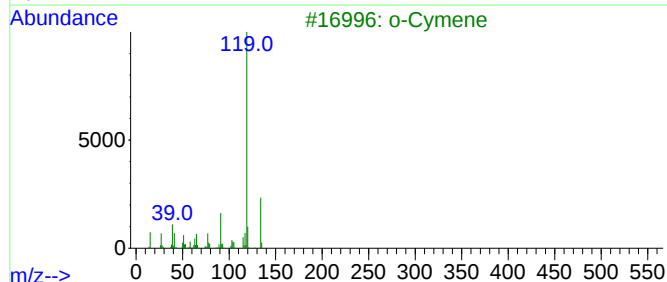
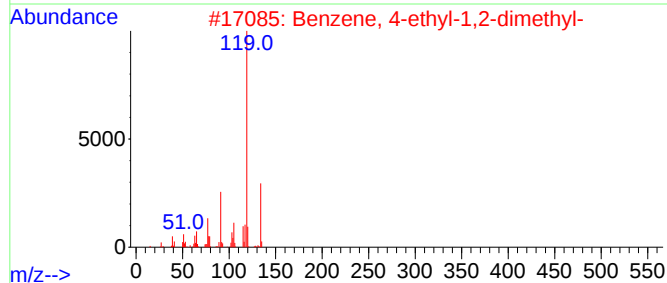
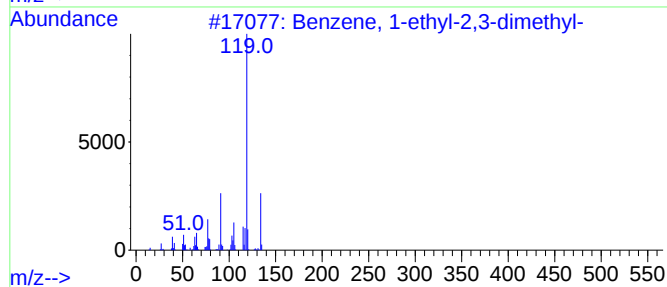
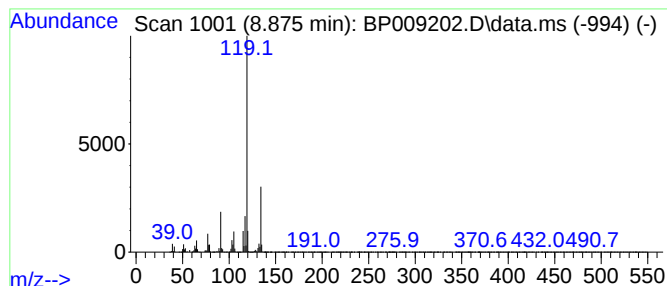
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.875	57.65 ng	3769140	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			o-Cymene	134	C10H14	000527-84-4	94
4			p-Cymene	134	C10H14	000099-87-6	93
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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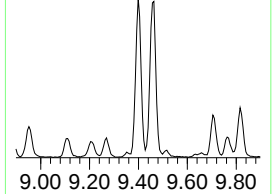
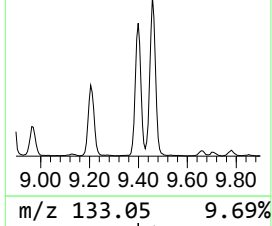
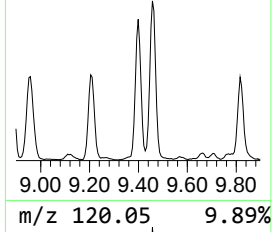
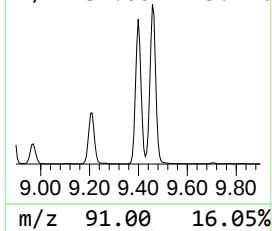
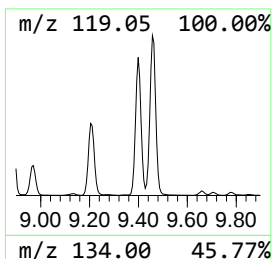
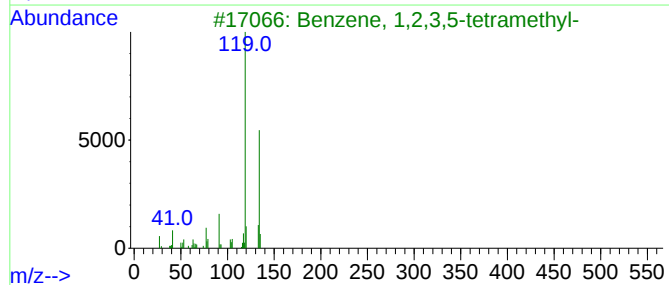
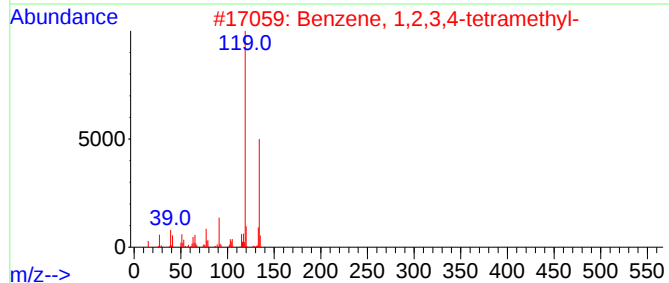
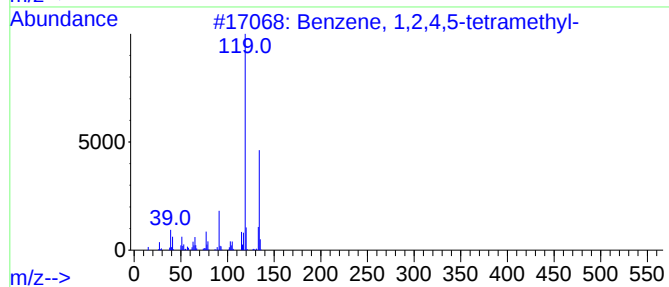
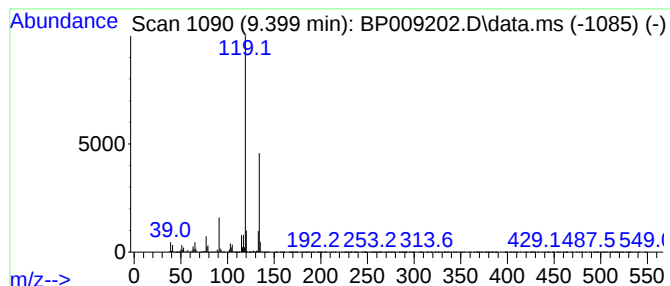
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.399	15.89 ng	1596850	Naphthalene-d8	10.569

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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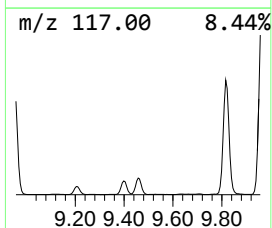
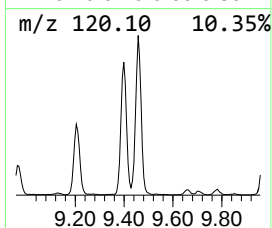
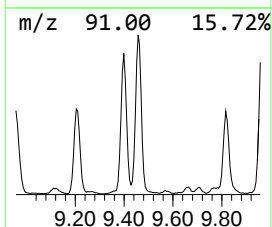
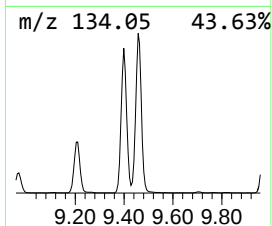
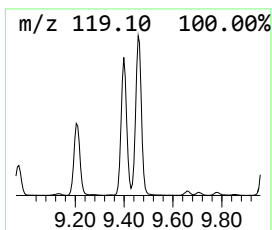
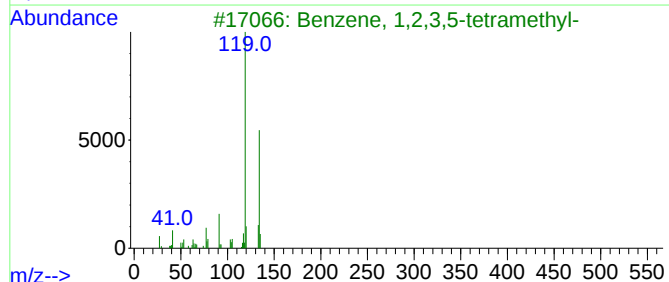
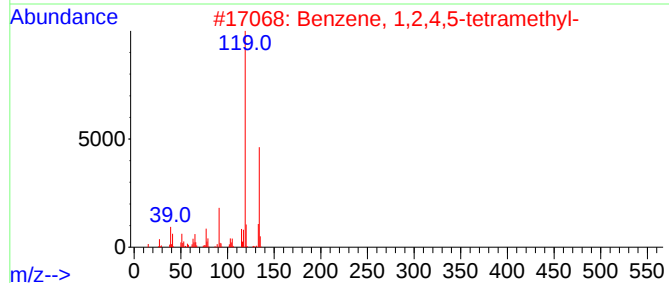
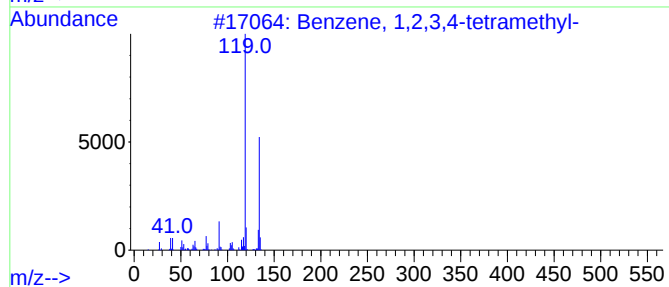
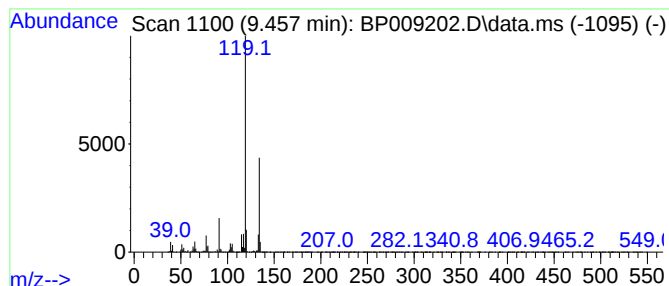
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.457	19.75 ng	1985250	Naphthalene-d8	10.569

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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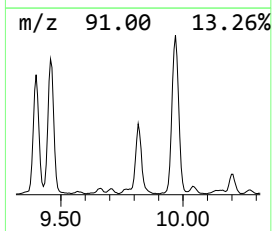
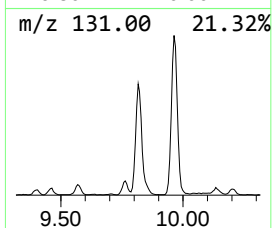
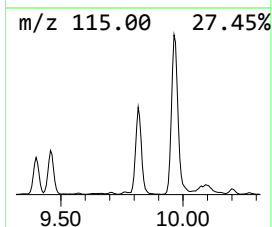
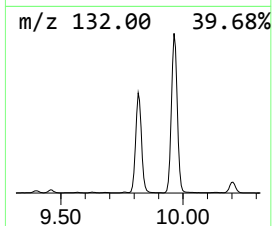
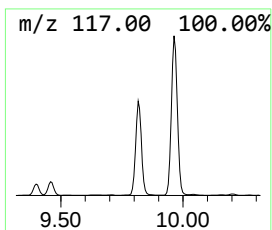
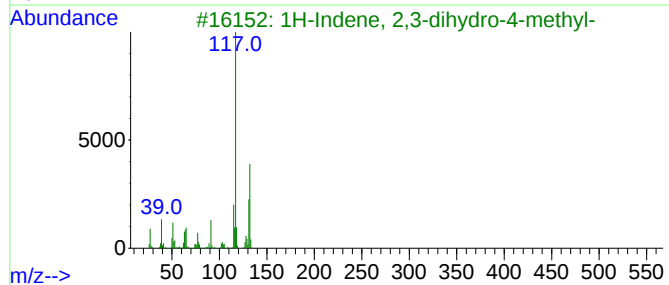
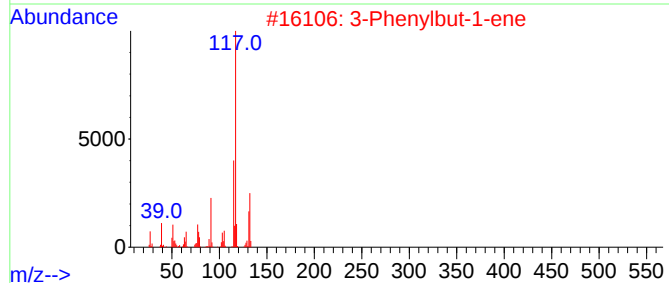
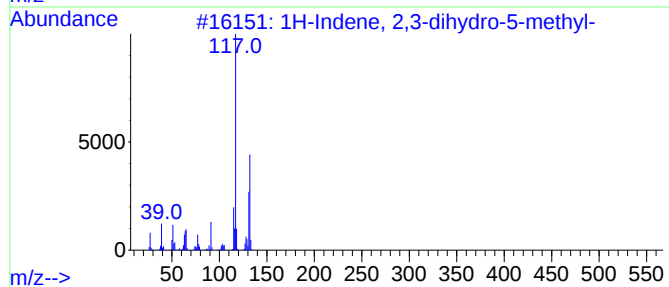
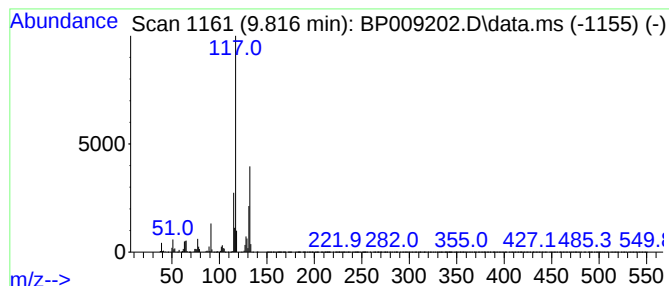
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.816	14.69 ng	1476130	Naphthalene-d8	10.569

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93	
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90	
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90	
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90	
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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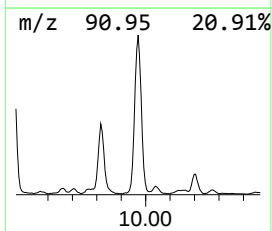
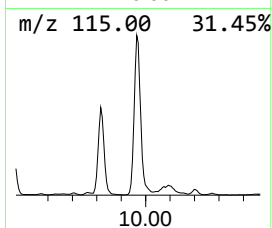
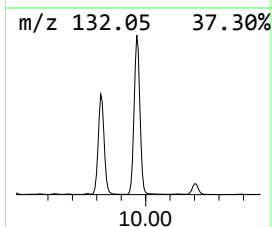
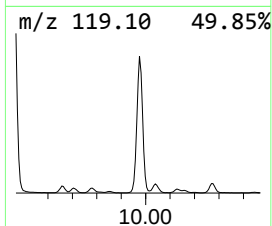
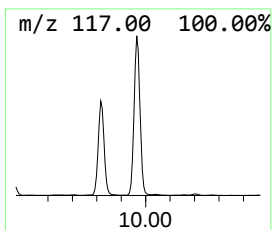
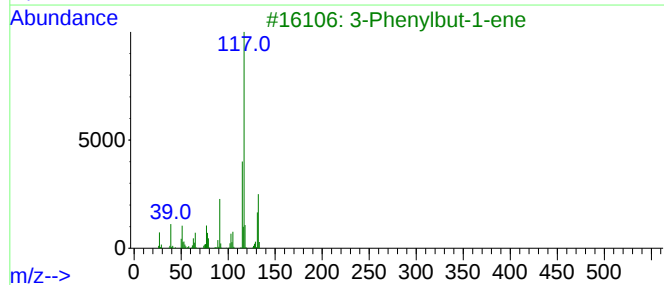
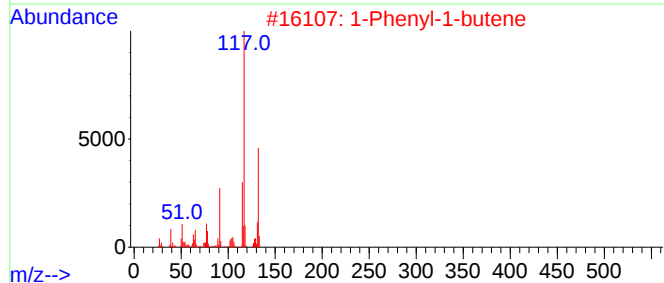
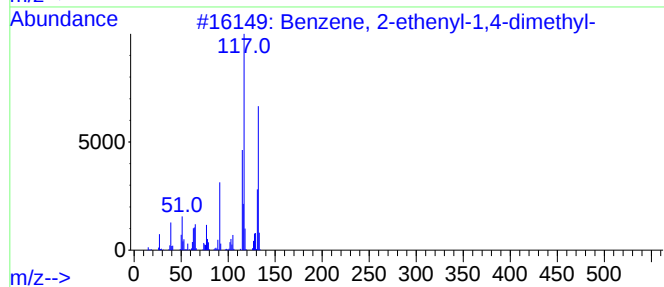
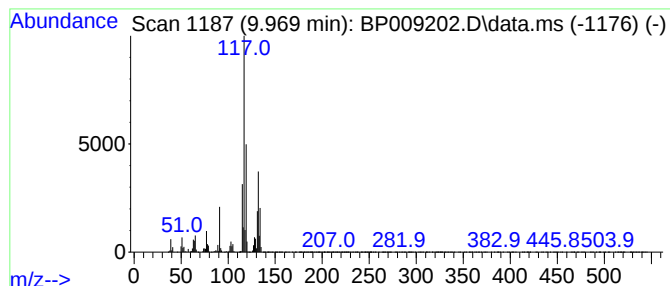
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.969	34.64 ng	3481500	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	70
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009202.D
Acq On : 17 Feb 2022 18:58
Operator : CG/JU
Sample : N1575-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

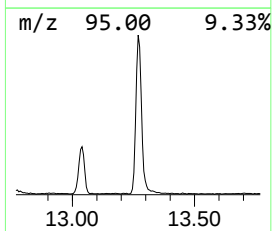
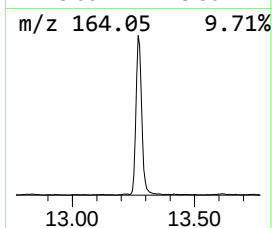
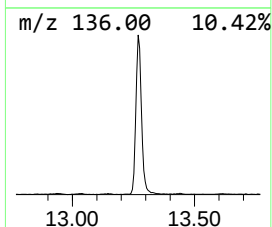
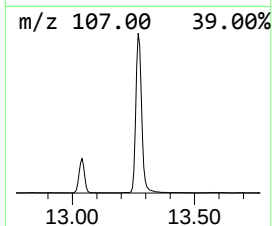
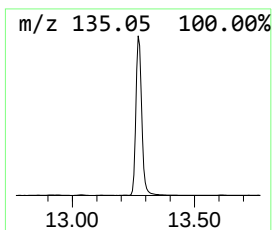
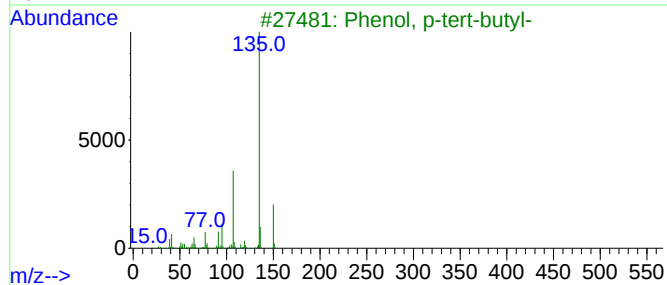
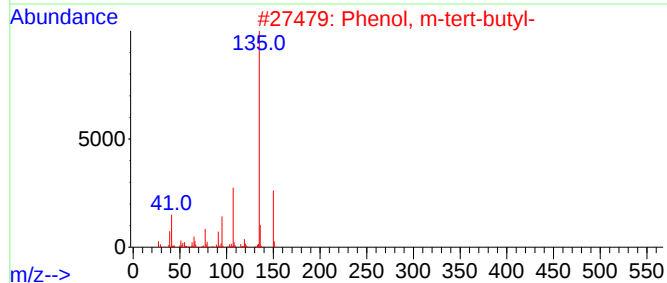
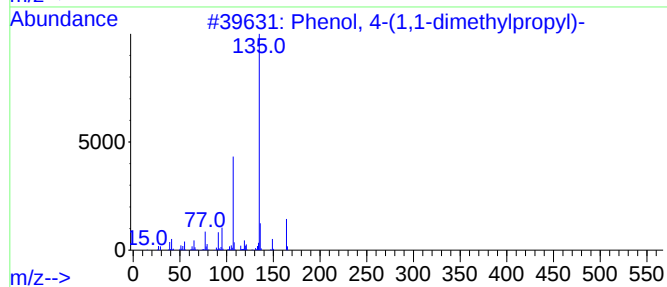
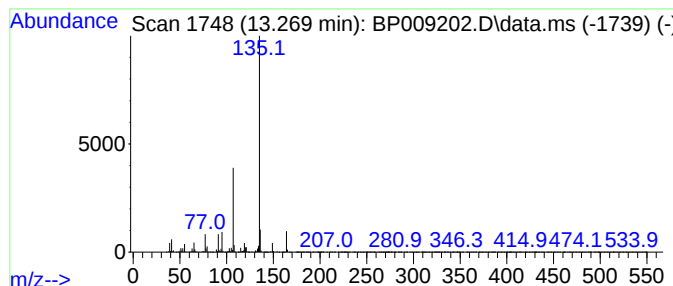
Instrument :
BNA_P
ClientSampleId :
MW-9R

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Phenol, 4-(1,1-dimethylprop... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.269	14.75 ng	2062690	Acenaphthene-d10			14.416
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-		164	C11H16O	000080-46-6	96
2	Phenol, m-tert-butyl-		150	C10H14O	000585-34-2	86
3	Phenol, p-tert-butyl-		150	C10H14O	000098-54-4	80
4	Hexestrol, O-trifluoroacetyl-		366	C20H21F3O3	1000365-31-7	64
5	4,6-Bis(4-ethoxybenzylthio)-5-ni...		457	C22H23N3O4S2	325957-76-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
 Data File : BP009202.D
 Acq On : 17 Feb 2022 18:58
 Operator : CG/JU
 Sample : N1575-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-9R

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.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
p-Xylene	5.787	74.6	ng	4876980	1	7.775	1307560	20.0
Benzene, (1-met...	6.258	26.2	ng	1710260	1	7.775	1307560	20.0
Benzene, propyl-	6.752	71.5	ng	4676690	1	7.775	1307560	20.0
Benzene, 1-ethy...	6.857	40.5	ng	2648460	1	7.775	1307560	20.0
Mesitylene	6.916	69.4	ng	4534750	1	7.775	1307560	20.0
Benzene, 1-ethy...	7.157	63.6	ng	4156110	1	7.775	1307560	20.0
Benzene, 1,2,3-...	7.422	316.0	ng	20657600	1	7.775	1307560	20.0
Benzene, 1,2,4-...	7.887	117.1	ng	7657090	1	7.775	1307560	20.0
Indane	8.140	46.3	ng	3026880	1	7.775	1307560	20.0
Benzene, 1,3-di...	8.263	22.9	ng	1497740	1	7.775	1307560	20.0
Benzene, 1,4-di...	8.410	34.5	ng	2257640	1	7.775	1307560	20.0
Benzene, 2-ethy...	8.722	23.7	ng	1547250	1	7.775	1307560	20.0
o-Cymene	8.775	16.9	ng	1107960	1	7.775	1307560	20.0
Benzene, 1-ethy...	8.875	57.6	ng	3769140	1	7.775	1307560	20.0
Benzene, 1,2,4,...	9.399	15.9	ng	1596850	2	10.569	2010380	20.0
Benzene, 1,2,3,...	9.457	19.8	ng	1985250	2	10.569	2010380	20.0
1H-Indene, 2,3-...	9.816	14.7	ng	1476130	2	10.569	2010380	20.0
Benzene, 2-ethe...	9.969	34.6	ng	3481500	2	10.569	2010380	20.0
Phenol, 4-(1,1-...	13.269	14.8	ng	2062690	3	14.416	2796400	20.0