

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
 Data File : BP003543.D
 Acq On : 07 Oct 2020 06:08
 Operator : CG/JU
 Sample : L4250-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-06

Integration Parameters: rteint.p

Integrator: RTE

Smoother: ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP003543.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.464	93	98	107	rVB	45028	72973	1.21%	0.255%
2	3.558	109	114	120	rBV	60976	94504	1.56%	0.330%
3	5.105	372	377	404	rVB	658525	1163115	19.26%	4.061%
4	6.705	637	649	676	rBV	355507	777654	12.88%	2.715%
5	7.493	776	783	794	rBV	318301	528323	8.75%	1.844%
6	7.857	839	845	857	rVB	160599	255769	4.23%	0.893%
7	7.987	859	867	872	rBV	115426	178281	2.95%	0.622%
8	8.134	886	892	896	rBV	46073	71947	1.19%	0.251%
9	8.181	896	900	908	rVB2	37100	61365	1.02%	0.214%
10	8.593	963	970	974	rBV	198401	317180	5.25%	1.107%
11	8.646	974	979	1002	rVB	1049328	1993190	33.00%	6.959%
12	9.116	1054	1059	1068	rVB	127125	202087	3.35%	0.706%
13	9.675	1145	1154	1164	rBV2	54185	110797	1.83%	0.387%
14	10.269	1245	1255	1261	rBV	416242	750542	12.43%	2.620%
15	10.328	1261	1265	1272	rVB3	40777	74251	1.23%	0.259%
16	10.898	1356	1362	1374	rBV2	38923	98179	1.63%	0.343%
17	12.169	1568	1578	1586	rBV	54515	114643	1.90%	0.400%
18	12.763	1672	1679	1690	rBV	2794583	4169960	69.04%	14.558%
19	13.616	1818	1824	1836	rBV	212874	321718	5.33%	1.123%
20	14.145	1908	1914	1921	rBV	681323	1025731	16.98%	3.581%
21	14.216	1921	1926	1936	rVB	187785	288063	4.77%	1.006%
22	14.563	1979	1985	1993	rVB	77601	125751	2.08%	0.439%
23	15.210	2089	2095	2102	rBV	68840	107732	1.78%	0.376%
24	15.651	2163	2170	2183	rBV	1986810	3192801	52.86%	11.147%
25	16.910	2378	2384	2388	rBV	832352	1250286	20.70%	4.365%
26	16.951	2388	2391	2398	rVB	120638	175541	2.91%	0.613%
27	17.051	2403	2408	2415	rVB	62726	102332	1.69%	0.357%
28	17.333	2450	2456	2468	rBV	68207	123166	2.04%	0.430%
29	17.839	2538	2542	2548	rBV3	66322	114475	1.90%	0.400%
30	17.980	2561	2566	2571	rBV	59582	82275	1.36%	0.287%
31	18.939	2725	2729	2736	rBV	246838	316810	5.25%	1.106%
32	19.292	2781	2789	2794	rBV	205774	309400	5.12%	1.080%
33	19.492	2818	2823	2831	rBV	4785670	6039855	100.00%	21.086%
34	20.739	3031	3035	3052	rBV	427973	619909	10.26%	2.164%
35	21.016	3076	3082	3087	rBV	1353963	1689025	27.96%	5.897%
36	23.180	3443	3450	3464	rVB	916984	1723689	28.54%	6.018%

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

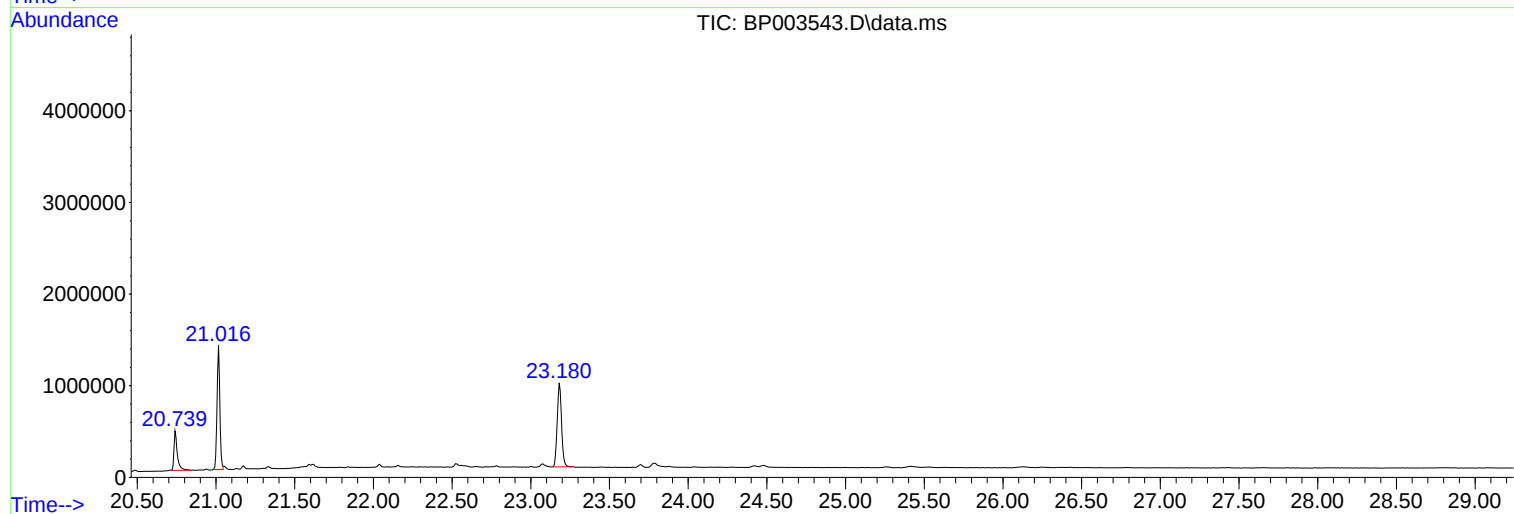
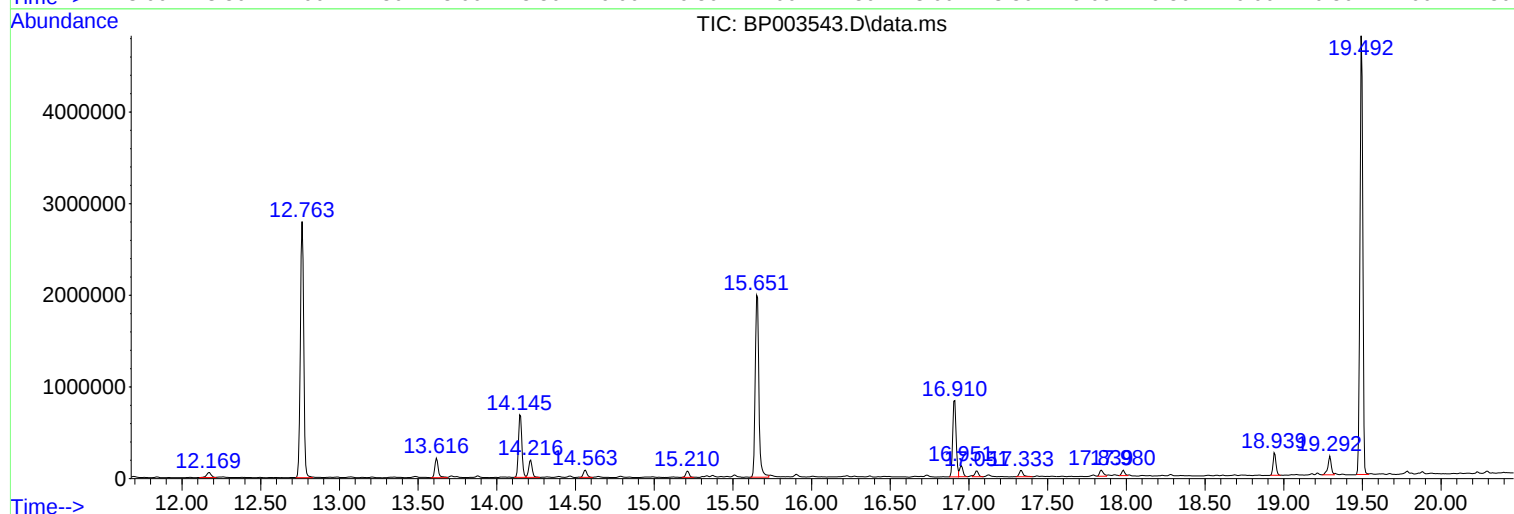
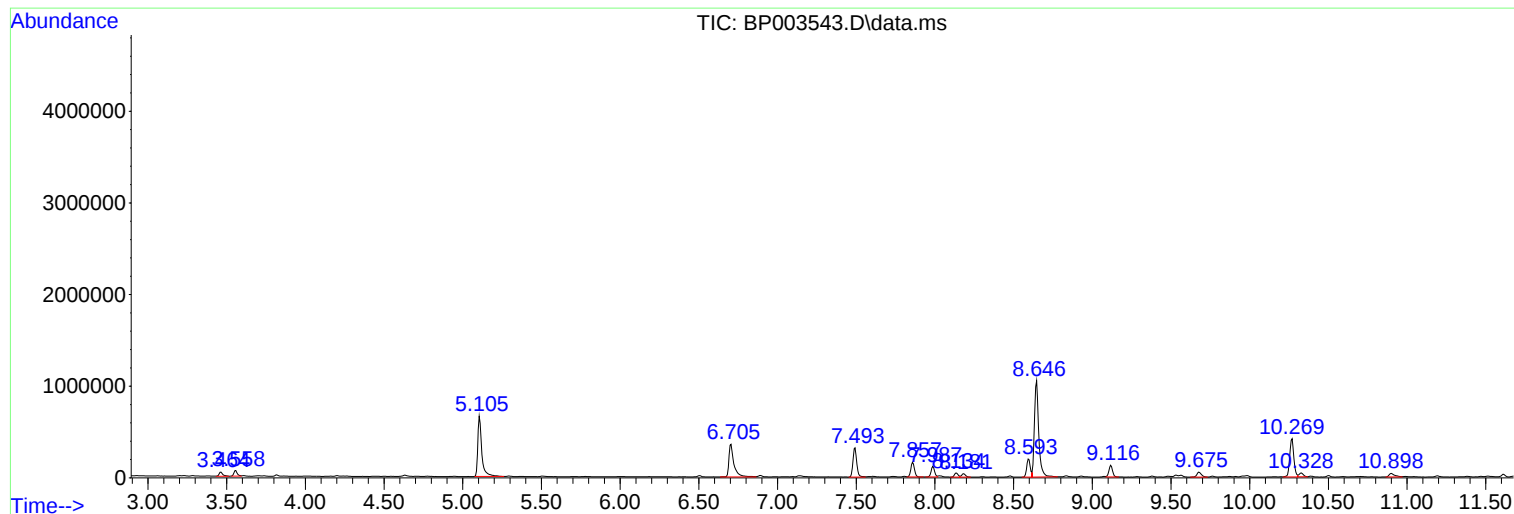
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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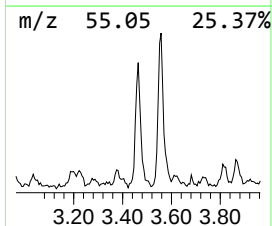
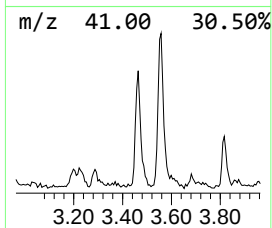
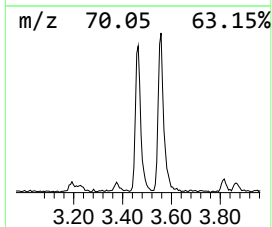
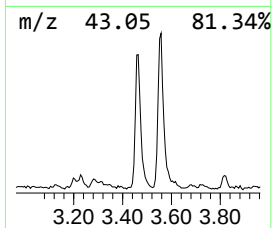
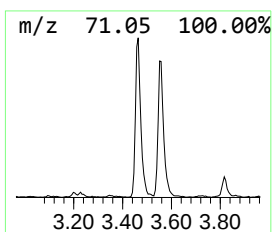
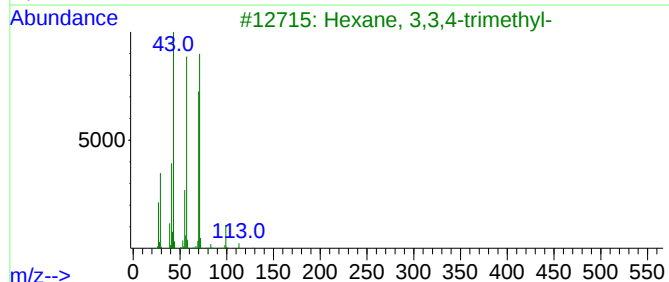
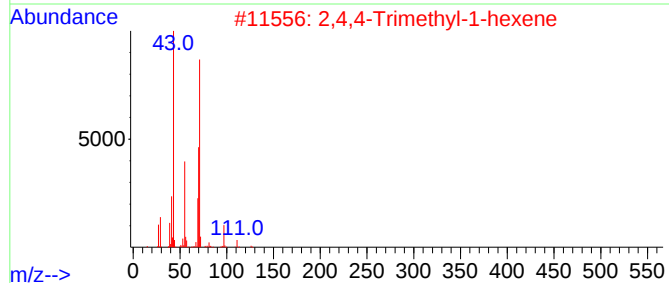
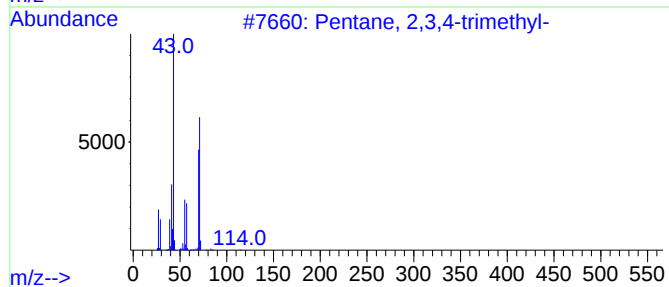
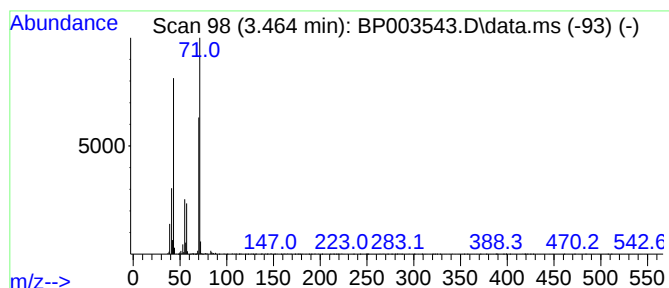
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Pentane, 2,3,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.464	2.76 ng	72973	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
3			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74
4			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
5			Pentane, 3-ethyl-	100	C7H16	000617-78-7	64



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

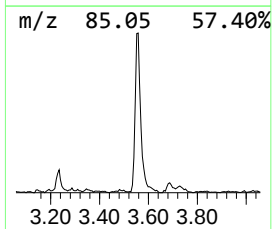
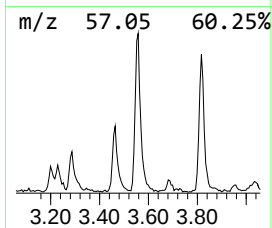
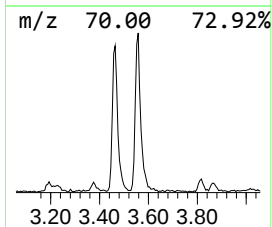
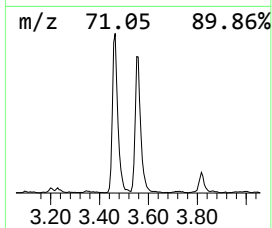
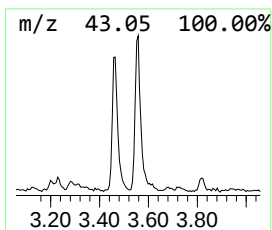
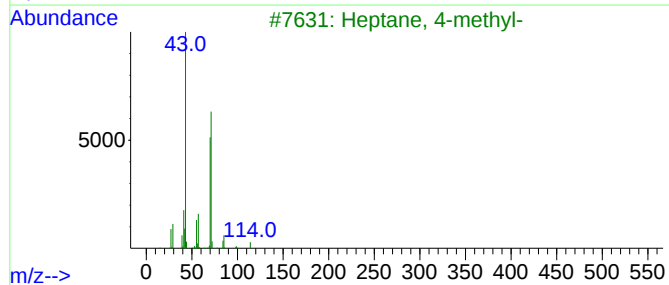
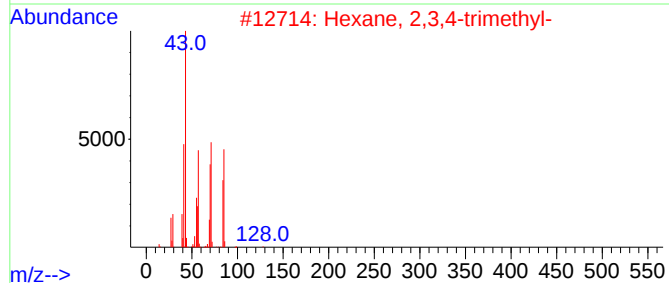
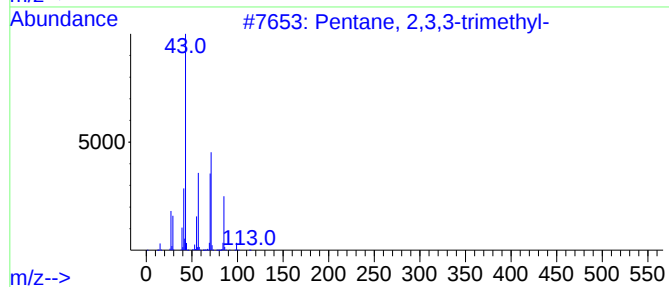
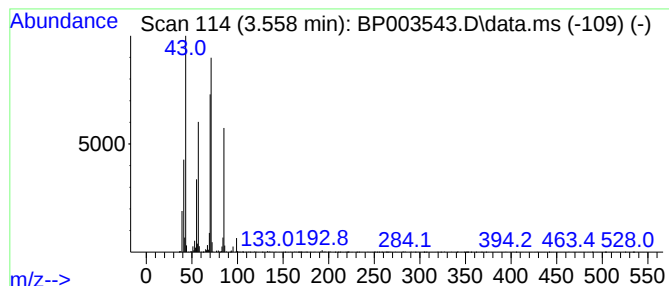
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.558	3.58 ng	94504	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
2			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	58
5			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	56



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

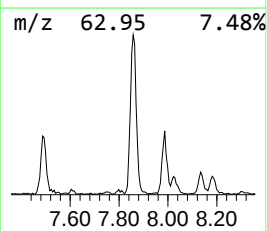
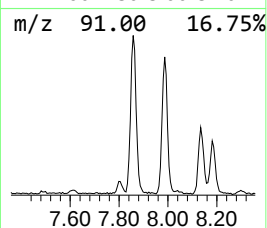
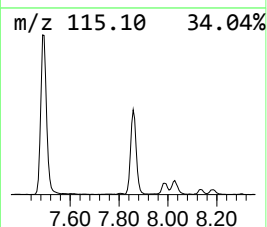
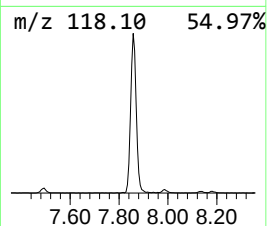
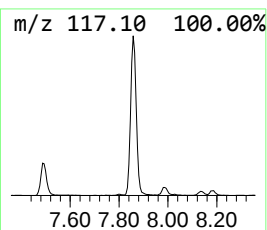
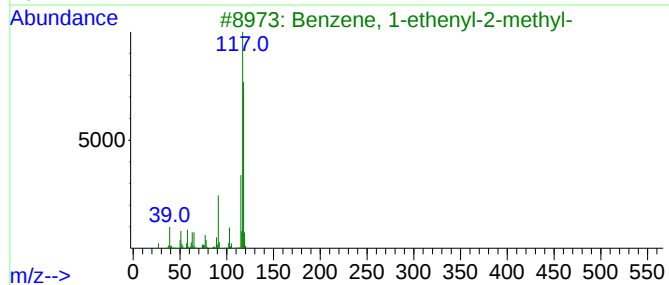
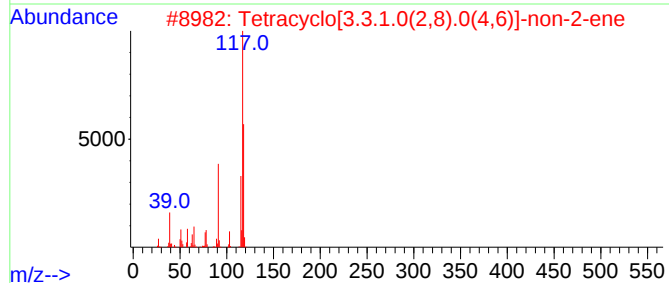
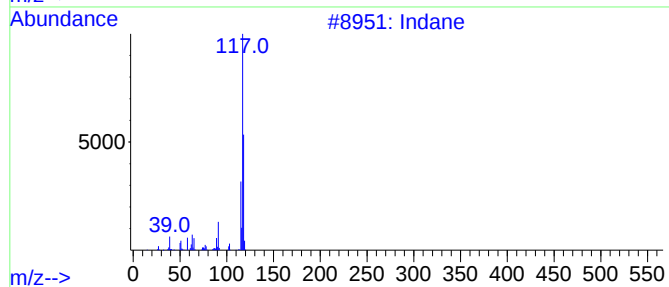
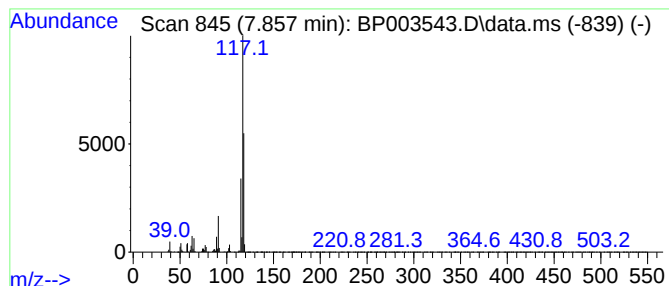
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.857	9.68 ng	255769	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	81
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	72
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	59



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ALS Vial : 22 Sample Multiplier: 1

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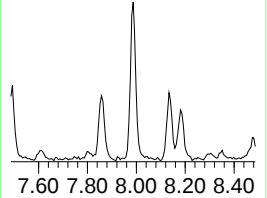
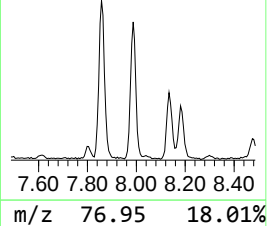
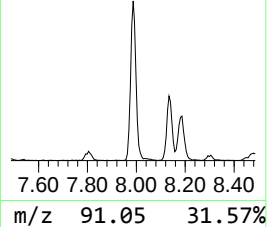
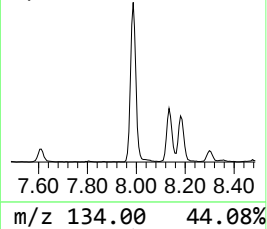
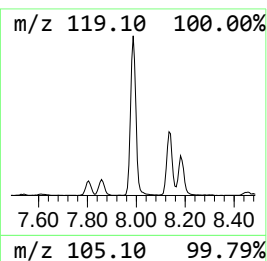
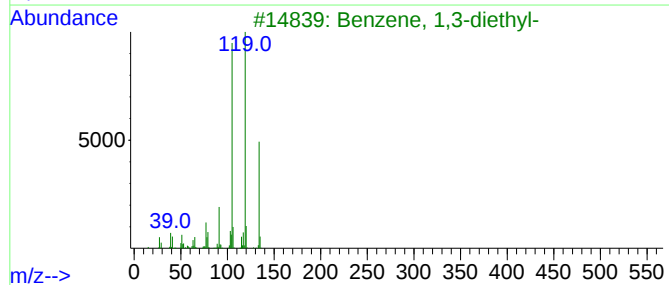
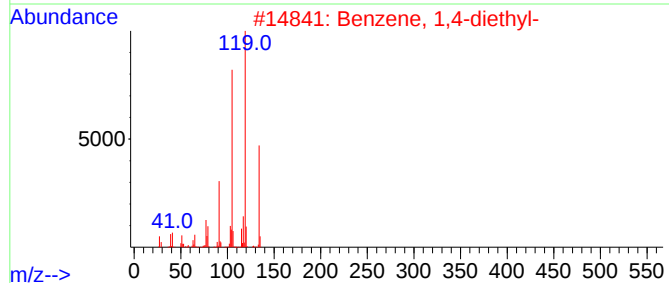
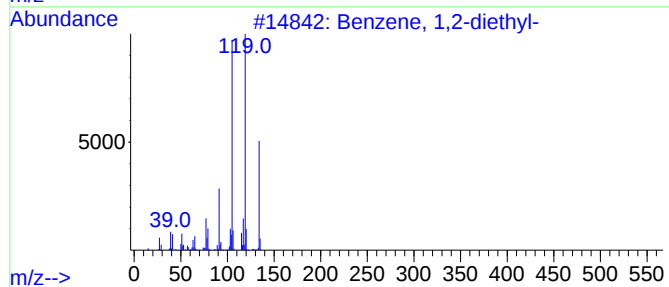
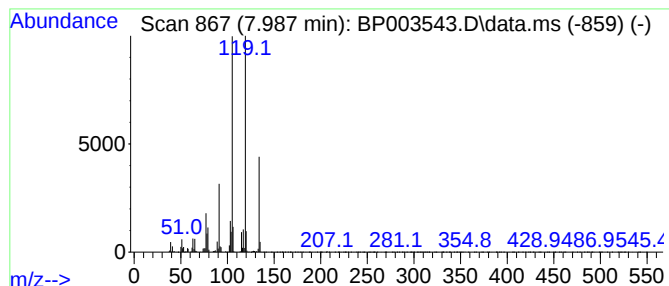
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	6.75 ng	178281	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



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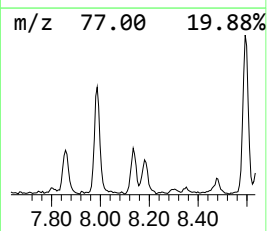
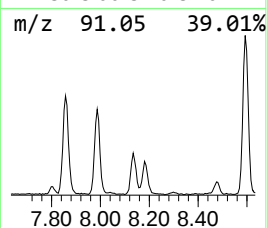
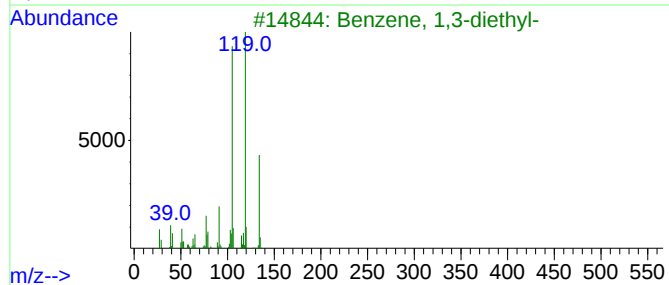
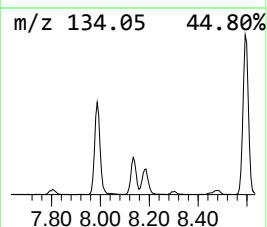
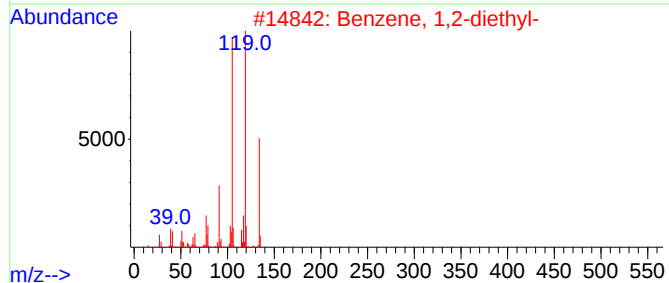
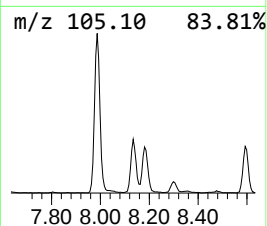
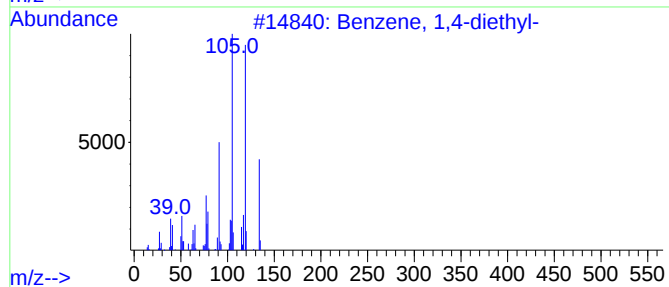
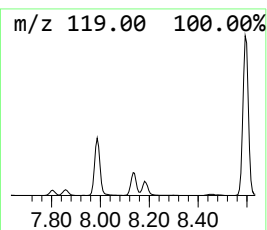
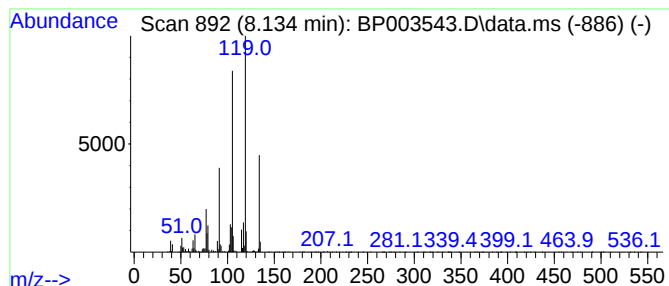
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.134	2.72 ng	71947	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	98
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

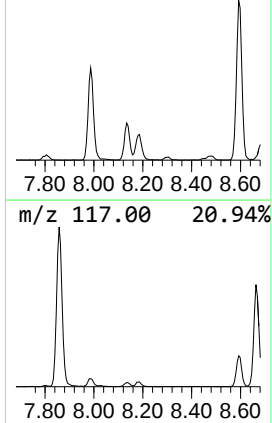
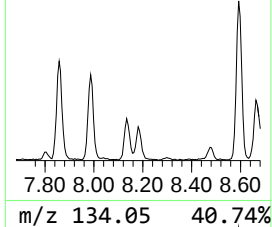
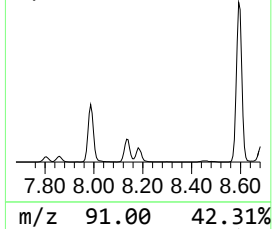
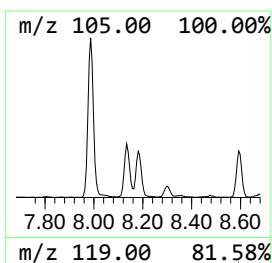
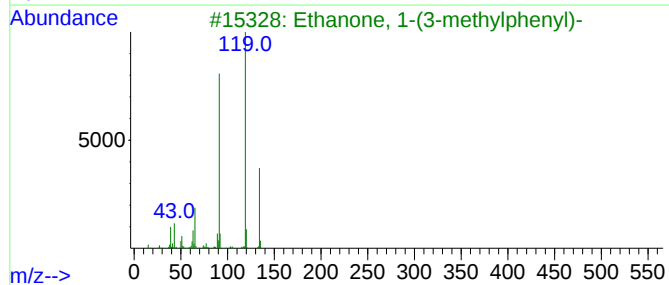
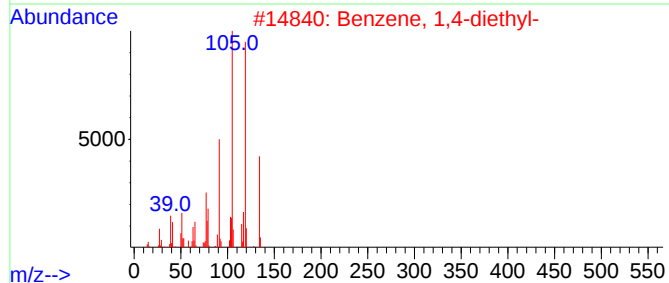
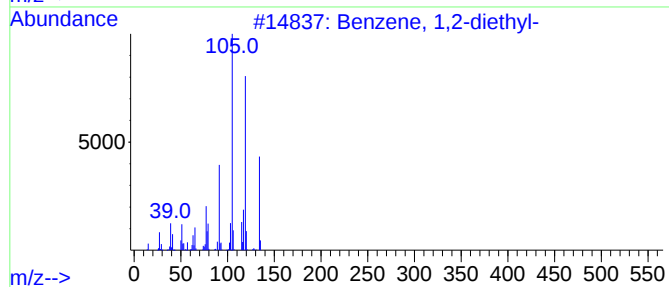
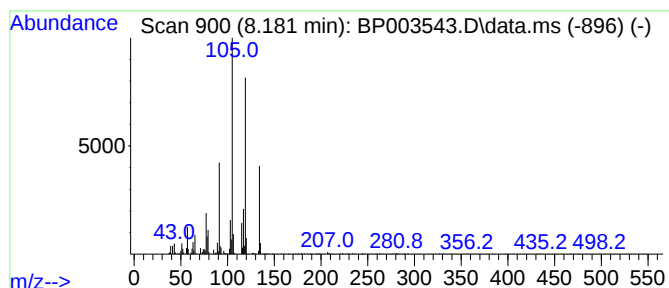
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanone, 1-(3-methylphenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	2.32 ng	61365	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3			Ethanone, 1-(3-methylphenyl)-	134	C9H10O	000585-74-0	91
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

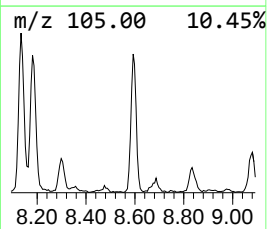
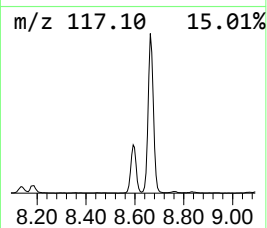
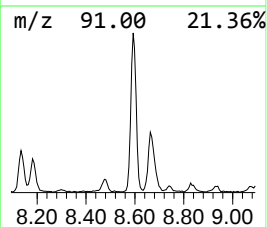
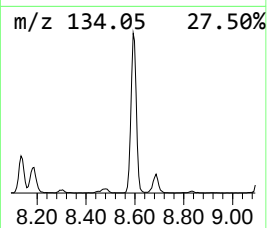
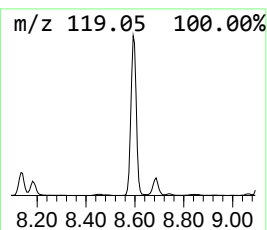
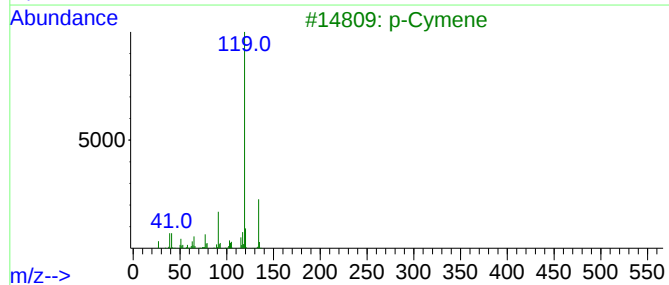
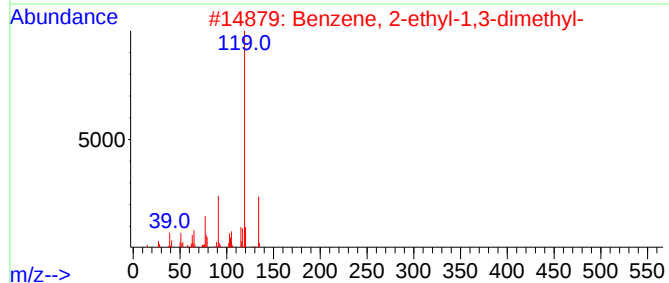
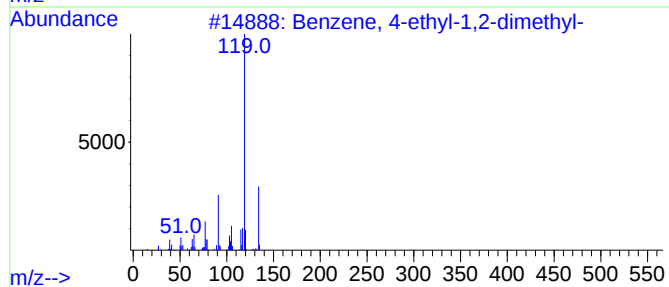
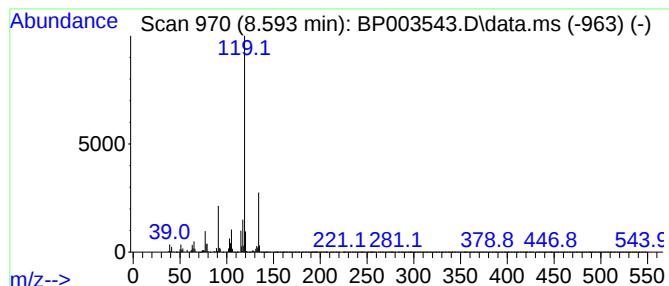
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.593	12.01 ng	317180	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

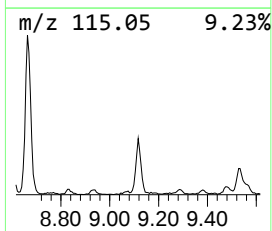
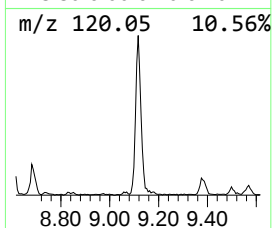
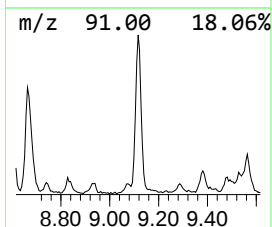
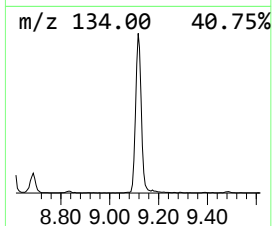
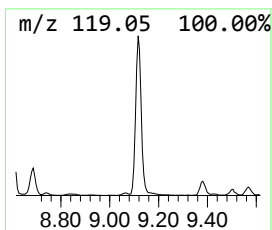
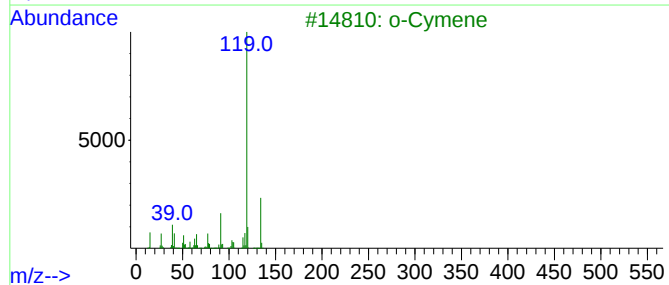
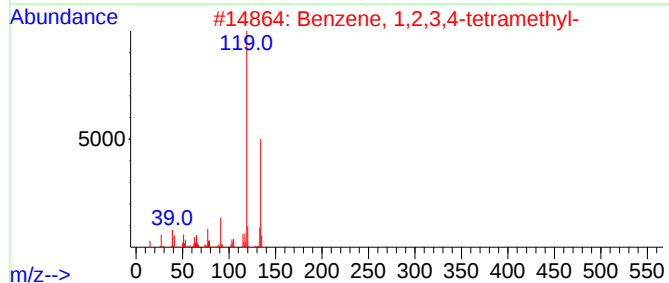
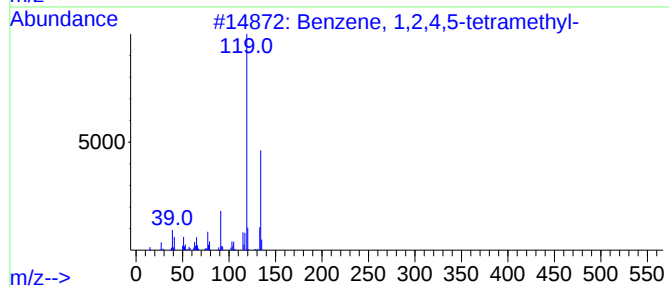
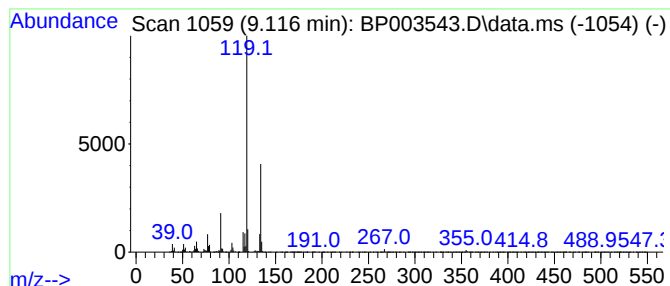
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.116	5.39 ng	202087	Naphthalene-d8	10.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			o-Cymene	134	C10H14	000527-84-4	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_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

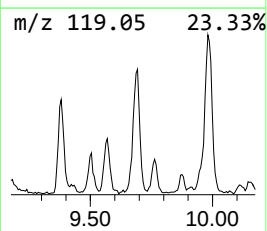
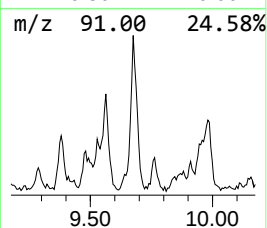
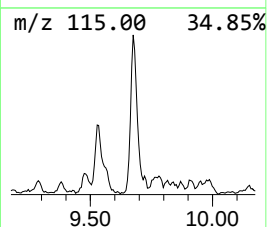
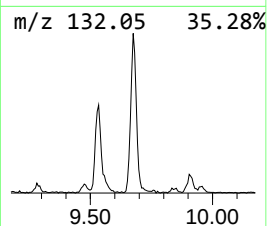
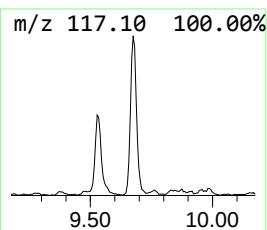
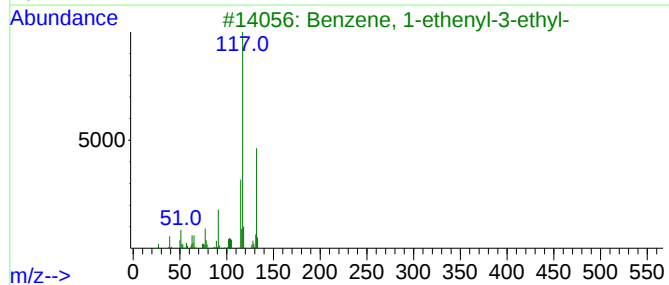
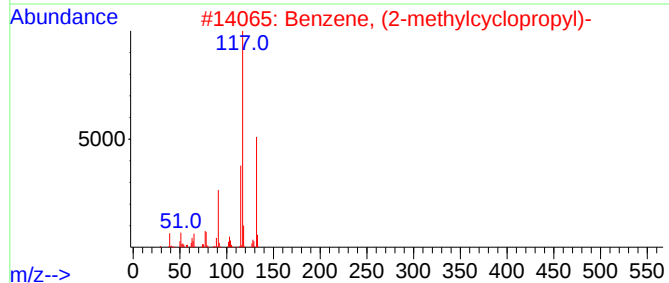
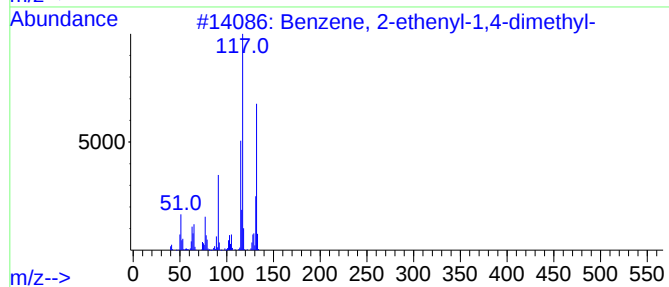
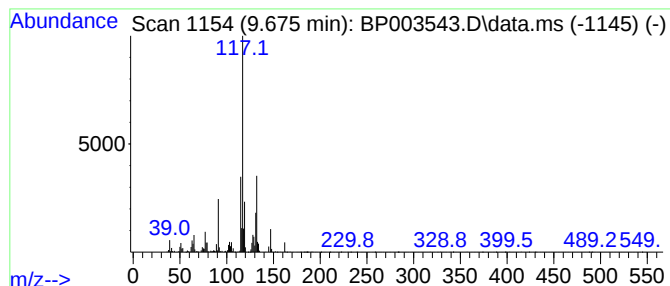
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.675	2.95 ng	110797	Naphthalene-d8	10.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	83
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

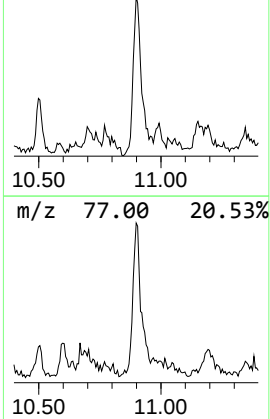
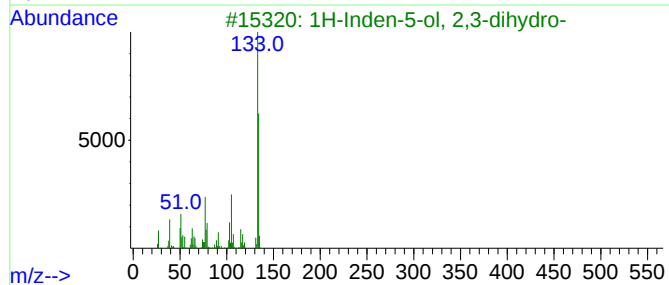
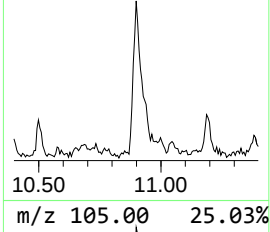
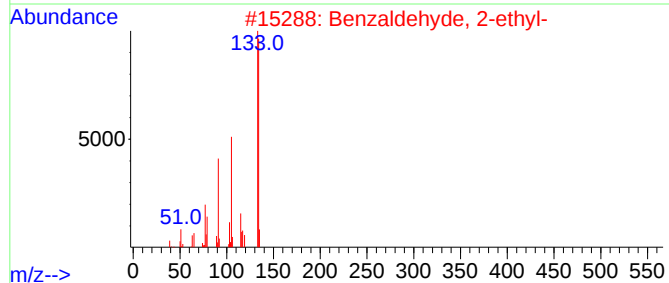
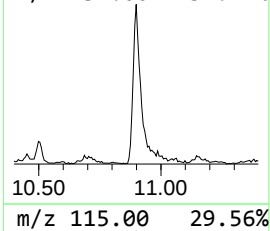
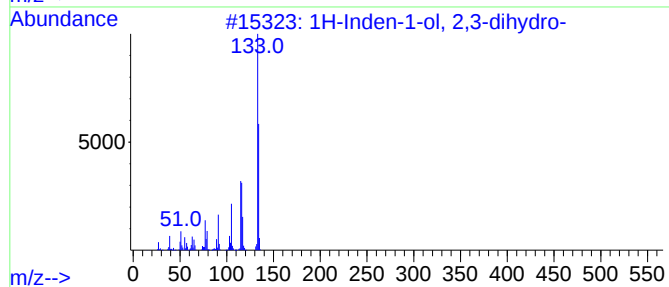
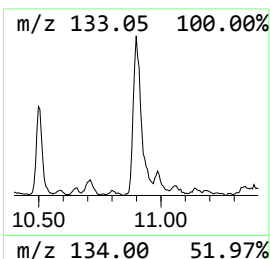
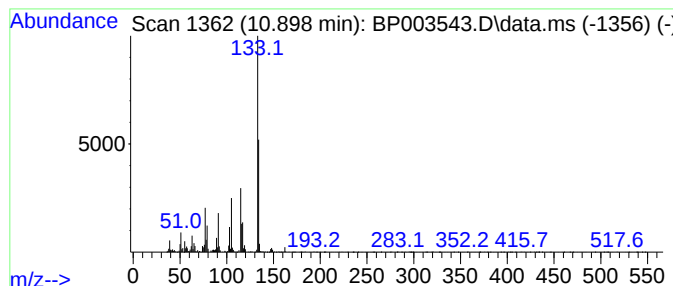
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1H-Inden-1-ol, 2,3-dihydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.898	2.62 ng	98179	Naphthalene-d8	10.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	87
2			Benzaldehyde, 2-ethyl-	134	C9H10O	022927-13-5	49
3			1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	49
4			1-(2,4-Dimethyl-phenyl)-2-(4H-[1...	247	C12H13N3OS	1000274-28-7	47
5			2,5-Pyrrolidinedione, 1-[(3,4-di...	247	C13H13N04	1000306-84-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

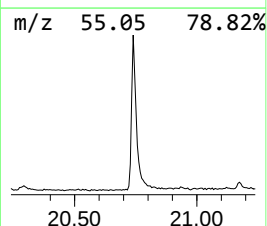
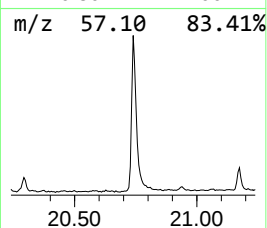
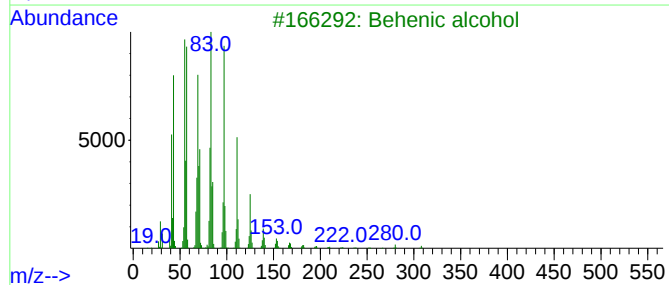
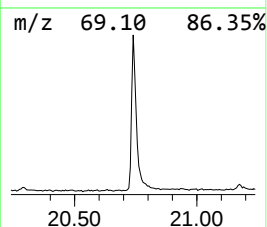
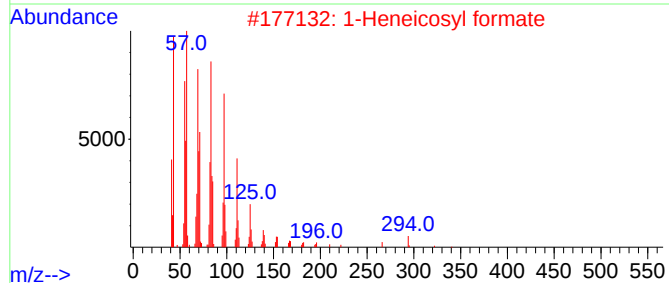
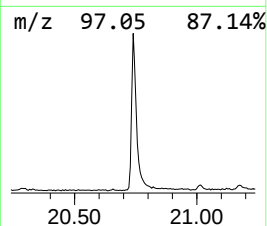
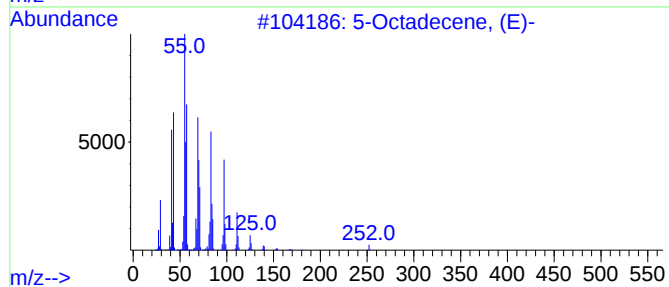
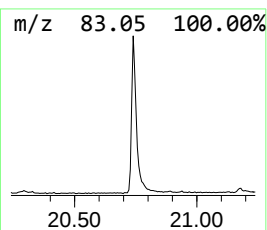
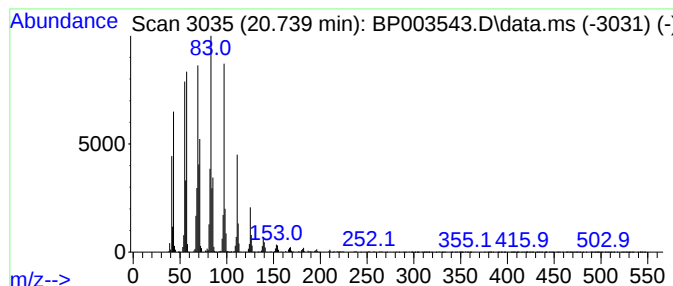
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 5-Octadecene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	7.34 ng	619909	Chrysene-d12	21.015

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	94
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3		Behenic alcohol	326	C22H46O	000661-19-8	94
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5		Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100620\
Data File : BP003543.D
Acq On : 07 Oct 2020 06:08
Operator : CG/JU
Sample : L4250-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-06

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,3,4-...	3.464	2.8	ng	72973	1	7.493	528323	20.0
Pentane, 2,3,3-...	3.558	3.6	ng	94504	1	7.493	528323	20.0
Indane	7.857	9.7	ng	255769	1	7.493	528323	20.0
Benzene, 1,2-di...	7.987	6.8	ng	178281	1	7.493	528323	20.0
Benzene, 1,4-di...	8.134	2.7	ng	71947	1	7.493	528323	20.0
Ethanone, 1-(3-...	8.181	2.3	ng	61365	1	7.493	528323	20.0
Benzene, 4-ethy...	8.593	12.0	ng	317180	1	7.493	528323	20.0
Benzene, 1,2,4,...	9.116	5.4	ng	202087	2	10.269	750542	20.0
Benzene, 2-ethe...	9.675	3.0	ng	110797	2	10.269	750542	20.0
1H-Inden-1-ol, ...	10.898	2.6	ng	98179	2	10.269	750542	20.0
5-Octadecene, (E)-	20.739	7.3	ng	619909	5	21.015	1689030	20.0