

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF011119\
 Data File : BF112034.D
 Acq On : 11 Jan 2019 23:49
 Operator : JU/SJ
 Sample : K1094-01 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CROSSWICKS-DRUM-1

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	206	213	220	rBV	185236	324978	1.20%	0.270%
2	5.369	552	558	561	rBV	6079900	6870573	25.32%	5.703%
3	5.498	570	580	585	rBV	10795677	27138883	100.00%	22.528%
4	5.751	615	623	625	rBV	9125119	12802252	47.17%	10.627%
5	6.045	670	673	677	rVB	375220	346613	1.28%	0.288%
6	6.345	719	724	728	rBV	1298836	1340143	4.94%	1.112%
7	6.422	730	737	739	rBV	5999588	7716026	28.43%	6.405%
8	6.451	739	742	744	rVV	4578686	3894979	14.35%	3.233%
9	6.498	744	750	753	rVB	5694636	5697187	20.99%	4.729%
10	6.581	758	764	767	rVB	5400037	5194450	19.14%	4.312%
11	6.740	781	791	794	rVV	11127483	22195716	81.79%	18.425%
12	6.822	802	805	806	rBV	378974	288516	1.06%	0.239%
13	6.840	806	808	811	rVB	413071	304208	1.12%	0.253%
14	6.916	817	821	823	rBV	774416	762742	2.81%	0.633%
15	6.951	823	827	830	rVB	5666284	6382330	23.52%	5.298%
16	7.063	843	846	849	rVV	2383964	1903755	7.01%	1.580%
17	7.128	853	857	859	rVV	1277234	1128216	4.16%	0.937%
18	7.157	859	862	865	rVV	2760671	2299770	8.47%	1.909%
19	7.204	865	870	872	rVV	4733720	3895744	14.35%	3.234%
20	7.275	878	882	885	rBV	985936	736553	2.71%	0.611%
21	7.345	890	894	896	rBV	1771577	1446434	5.33%	1.201%
22	7.369	896	898	901	rVV	1540395	1074885	3.96%	0.892%
23	7.416	902	906	908	rVV	2413863	1884081	6.94%	1.564%
24	7.645	942	945	948	rBV	549503	487275	1.80%	0.404%
25	7.675	948	950	953	rVB	784002	543308	2.00%	0.451%
26	10.833	1484	1487	1490	rBV	522307	502130	1.85%	0.417%
27	11.039	1519	1522	1526	rBV2	264915	390782	1.44%	0.324%
28	11.186	1543	1547	1548	rBV2	383417	508350	1.87%	0.422%
29	11.204	1548	1550	1552	rVV2	290446	329627	1.21%	0.274%
30	11.298	1563	1566	1574	rVB	903620	893047	3.29%	0.741%
31	11.998	1683	1685	1687	rBV2	436033	505940	1.86%	0.420%
32	12.769	1814	1816	1825	rBV5	251670	678123	2.50%	0.563%

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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_F\METHODS\8270-BF010719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

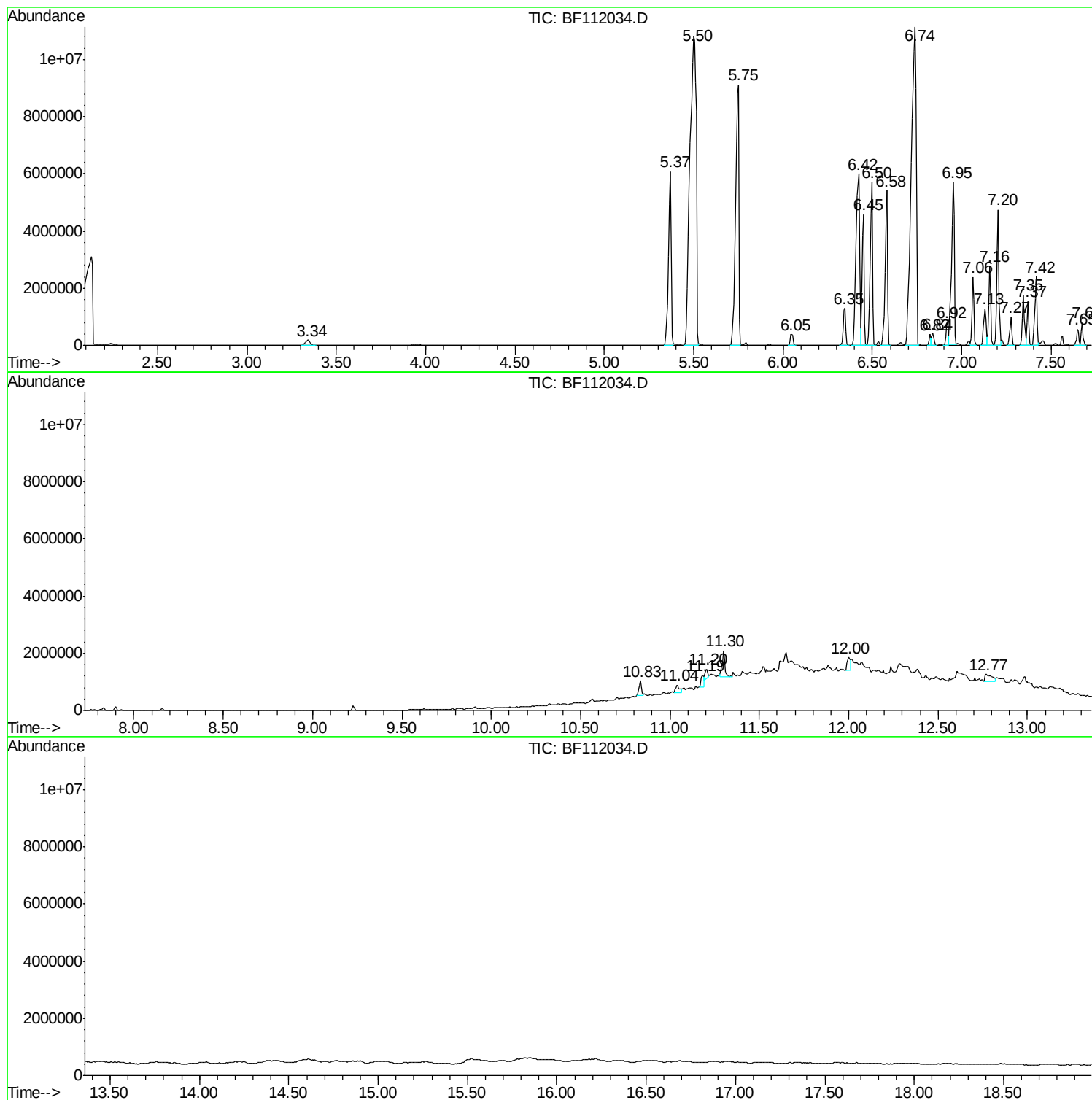
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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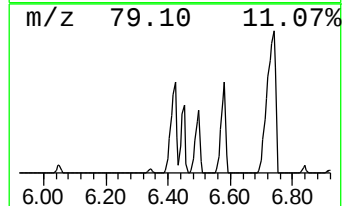
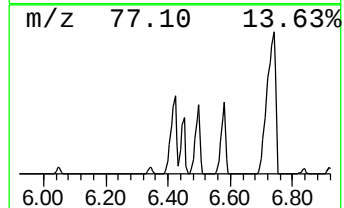
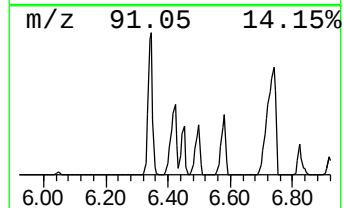
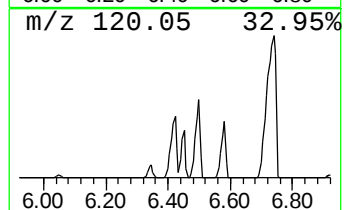
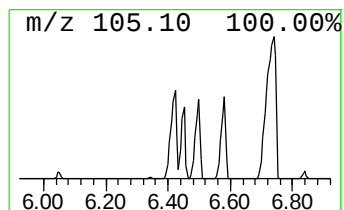
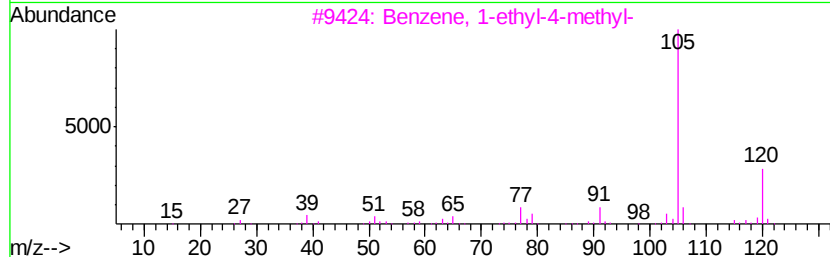
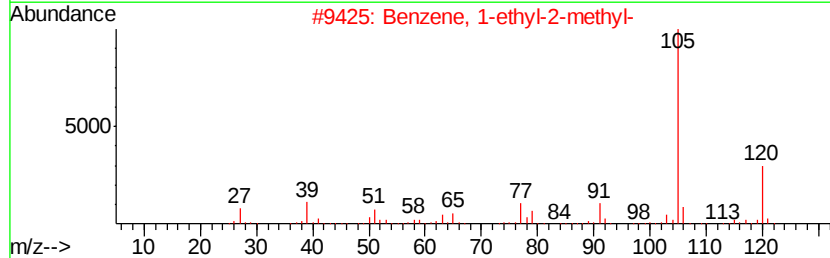
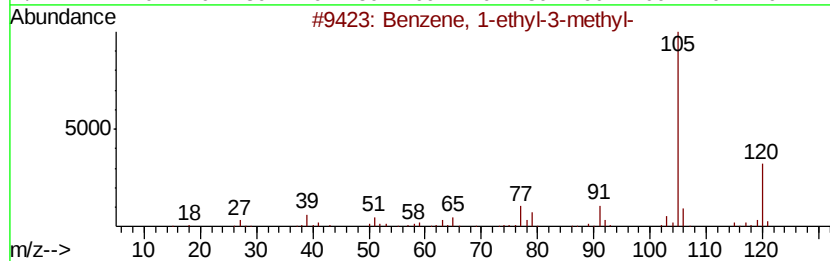
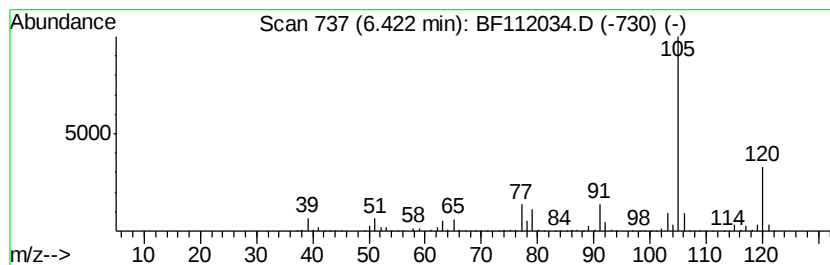
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	202.32 ng	7716030	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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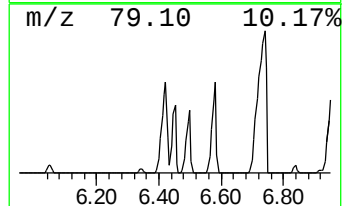
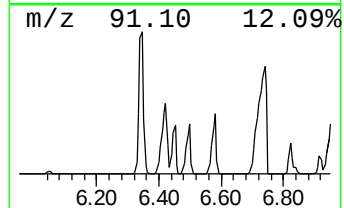
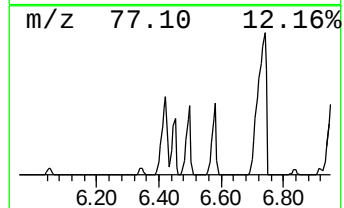
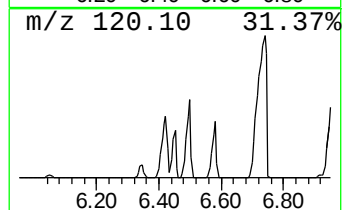
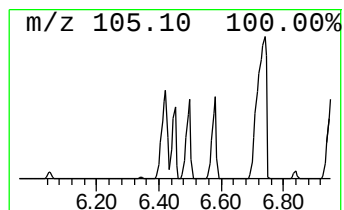
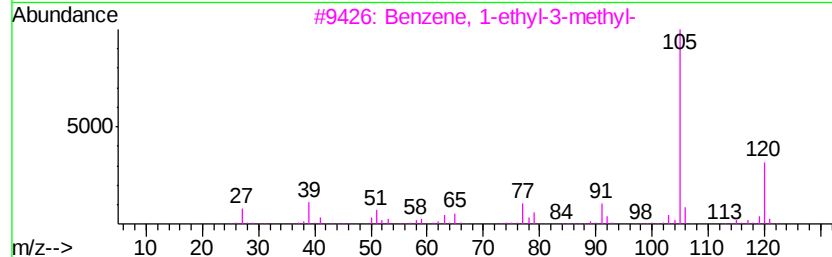
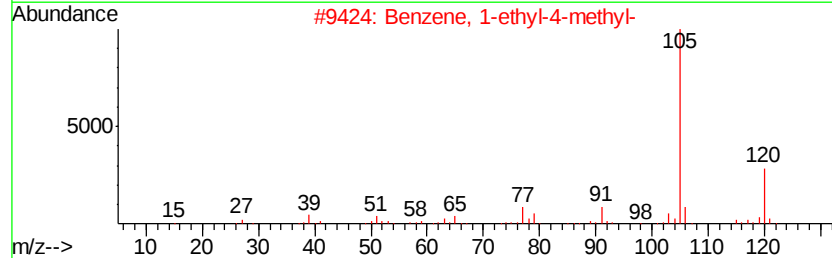
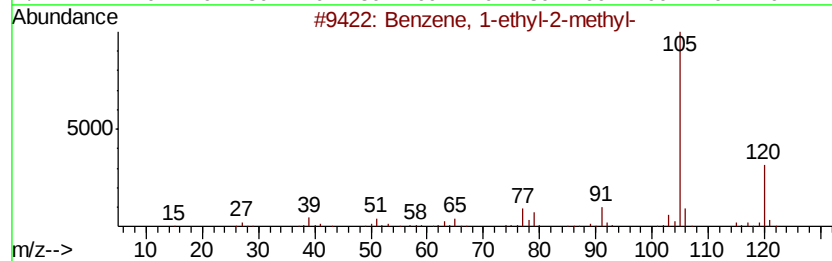
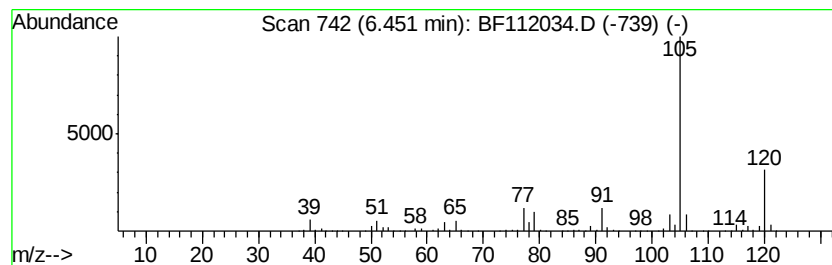
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	102.13 ng	3894980	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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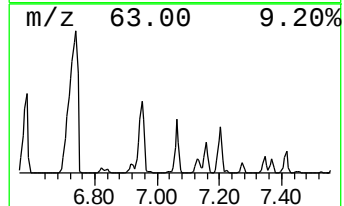
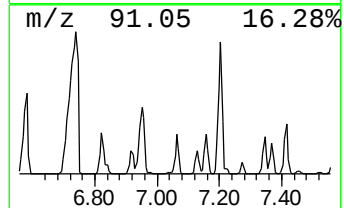
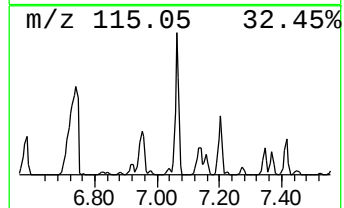
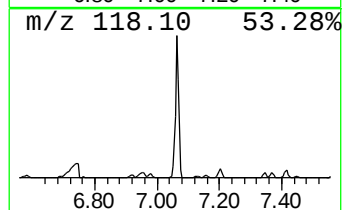
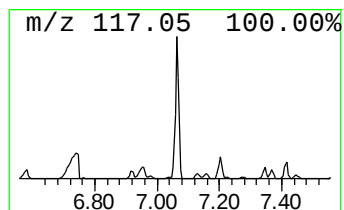
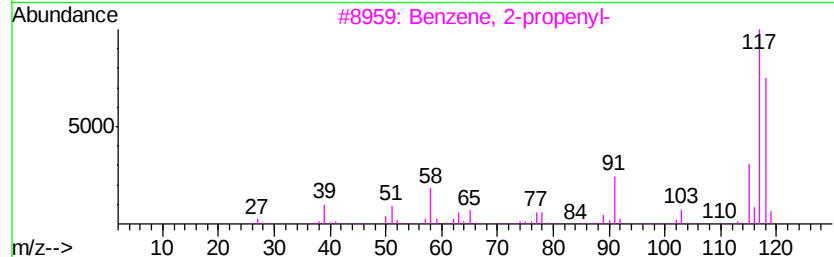
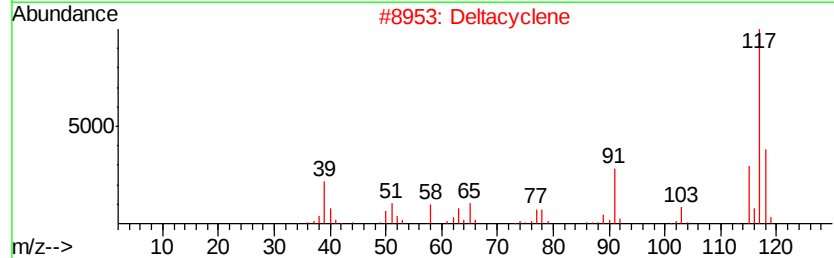
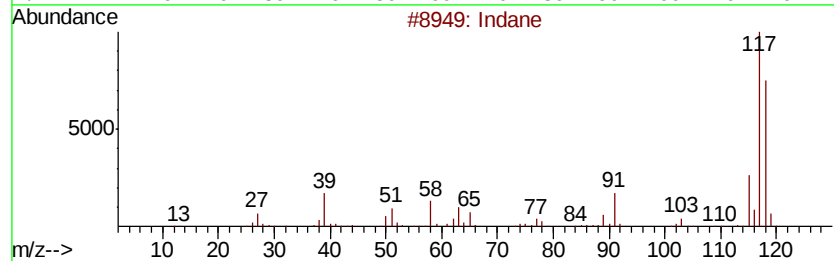
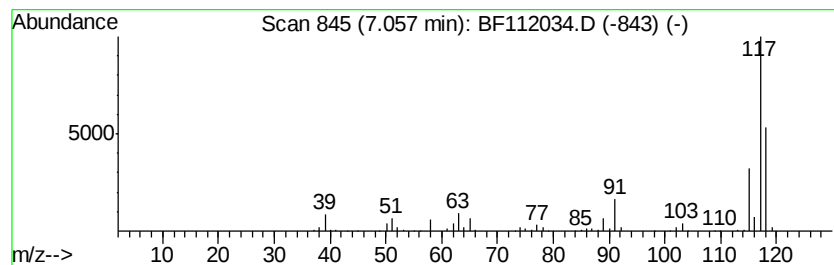
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	49.92 ng	1903760	1,4-Dichlorobenzene-d4	6.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Deltacyclene	118	C9H10	007785-10-6	72
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
5	Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	53



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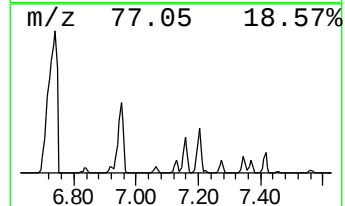
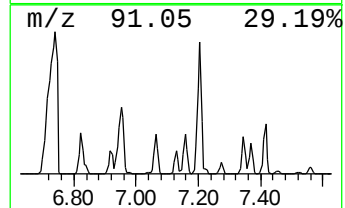
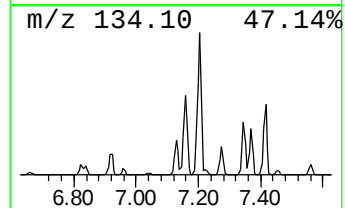
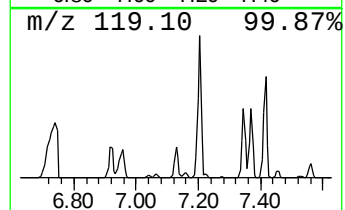
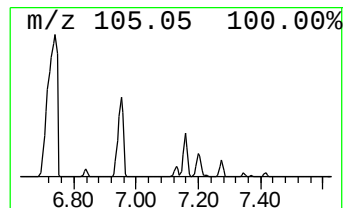
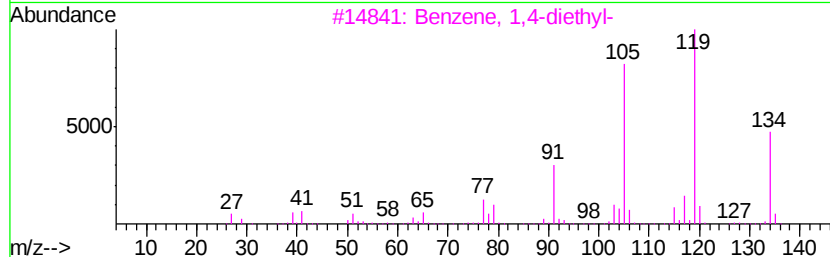
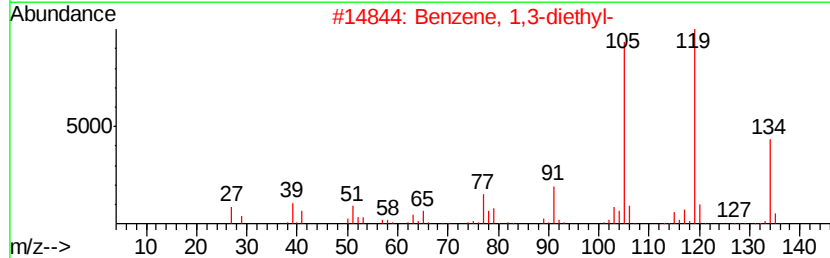
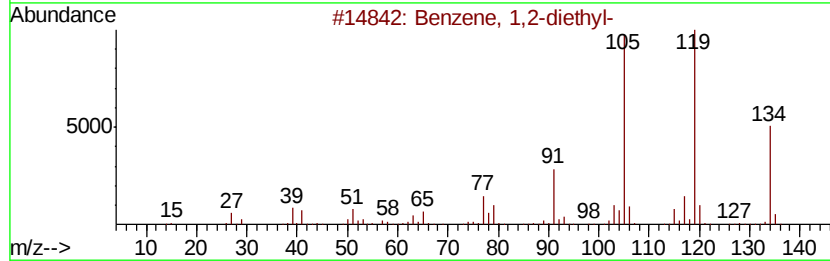
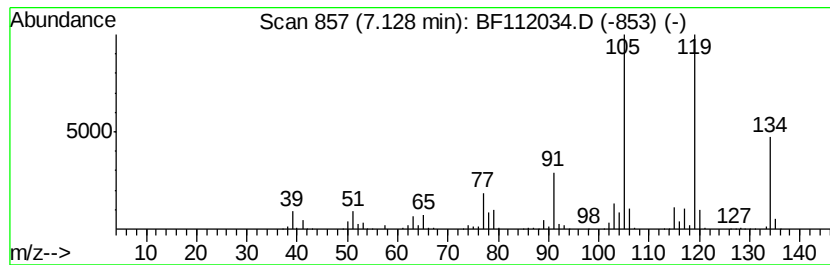
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,2-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	29.58 ng	1128220	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
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



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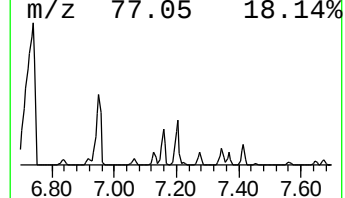
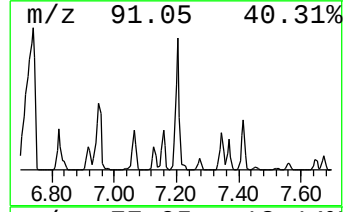
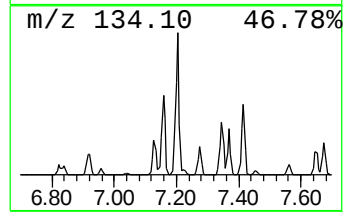
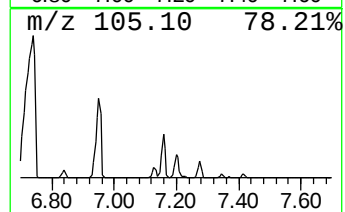
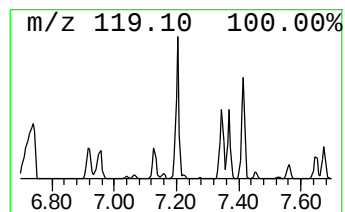
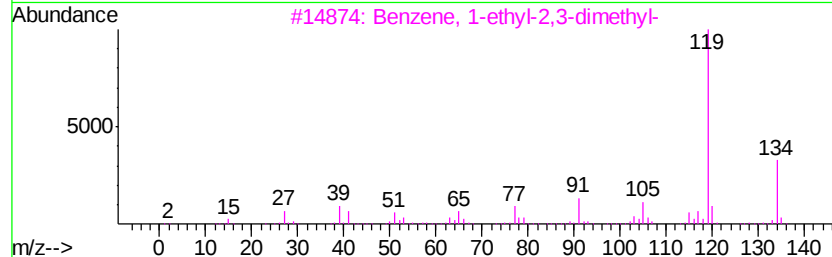
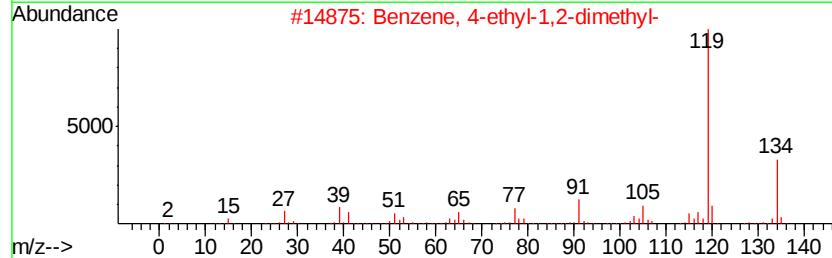
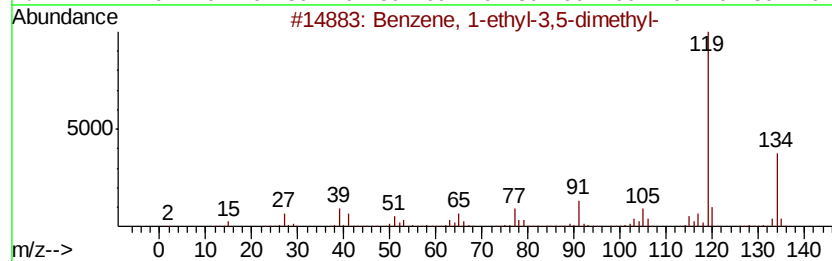
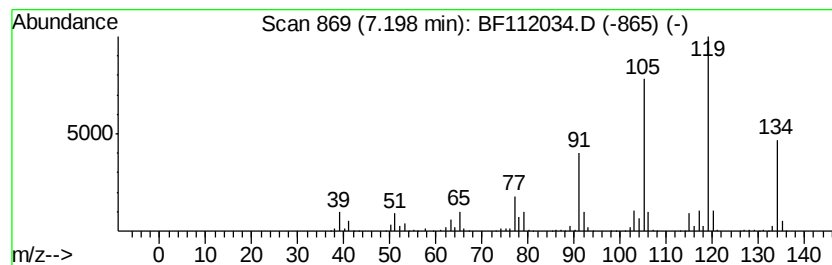
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Peak Number 11 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	102.15 ng	3895740	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
4		o-Cymene	134	C10H14	000527-84-4	93
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91



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Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-1

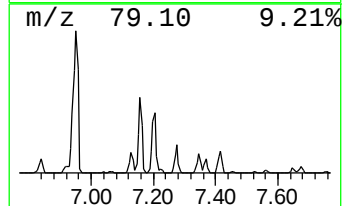
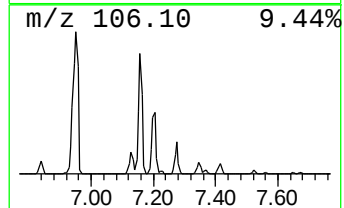
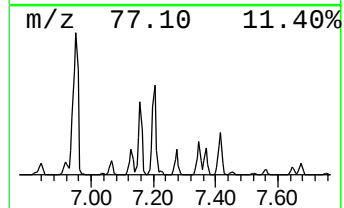
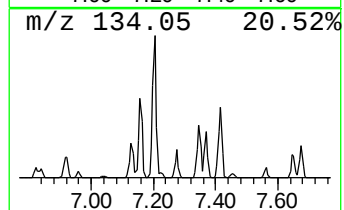
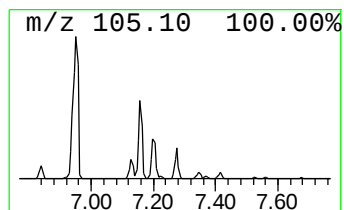
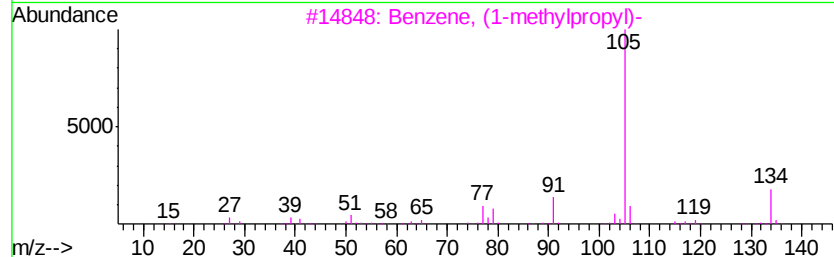
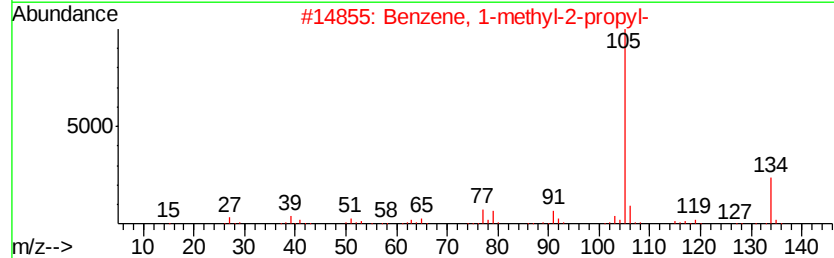
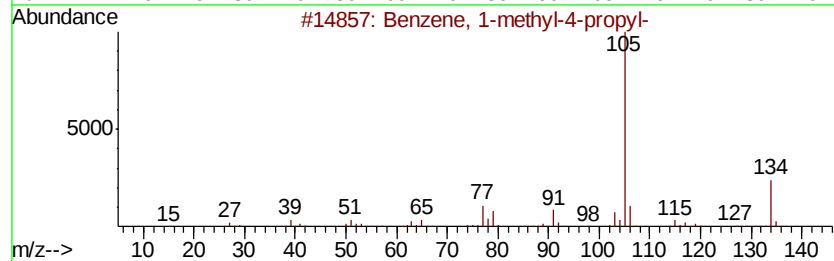
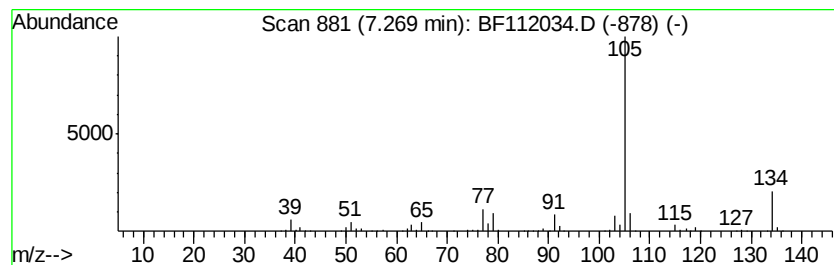
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1-methyl-4-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	19.31 ng	736553	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	90
5		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
Acq On : 11 Jan 2019 23:49
Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CROSSWICKS-DRUM-1

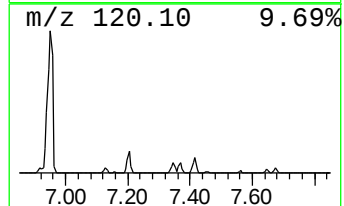
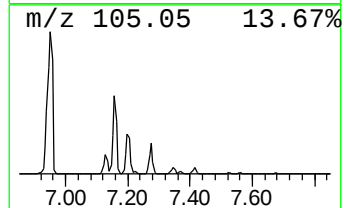
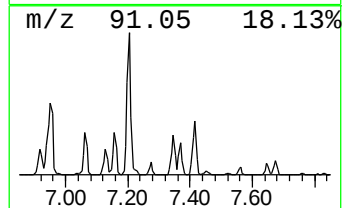
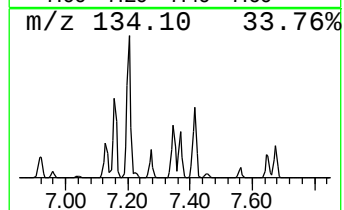
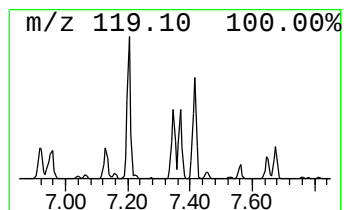
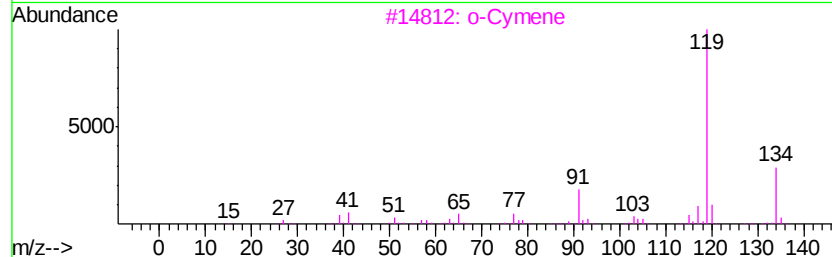
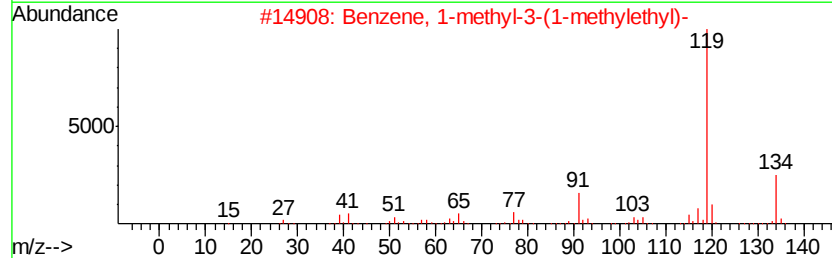
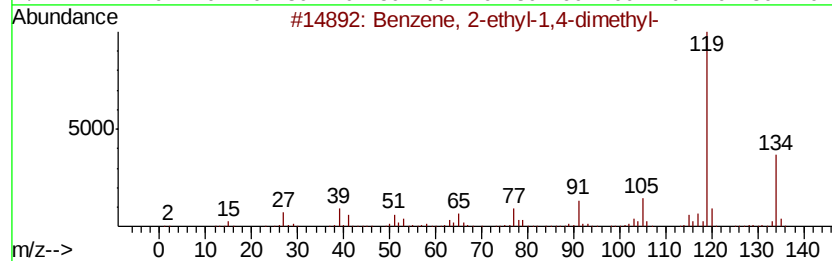
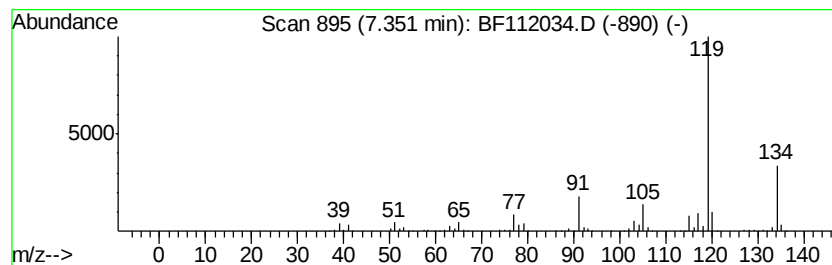
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.35	37.93 ng	1446430	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		p-Cymene	134	C10H14	000099-87-6	94
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
Acq On : 11 Jan 2019 23:49
Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CROSSWICKS-DRUM-1

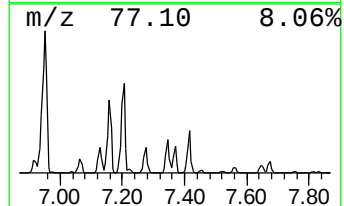
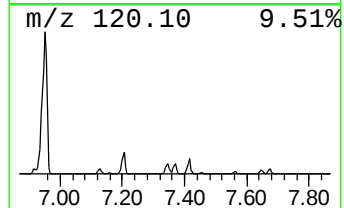
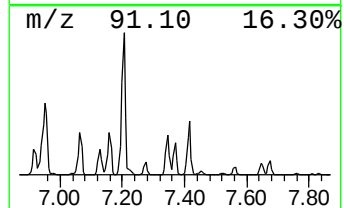
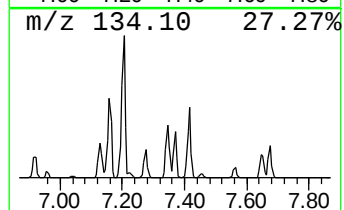
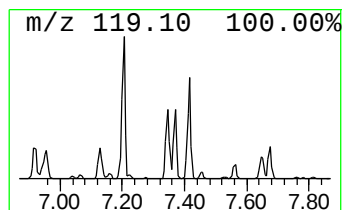
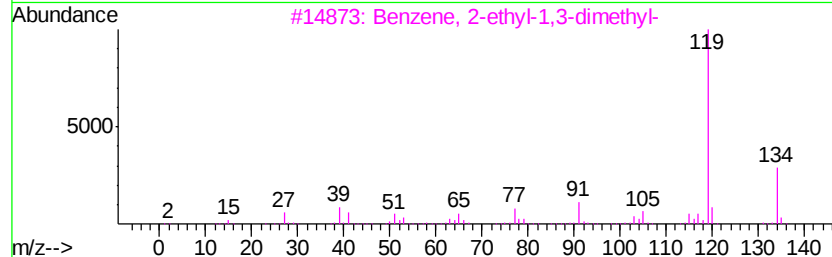
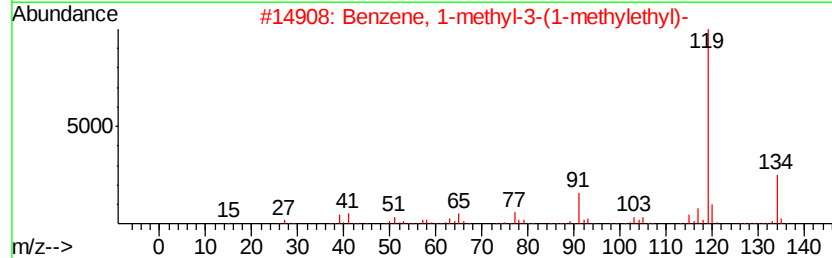
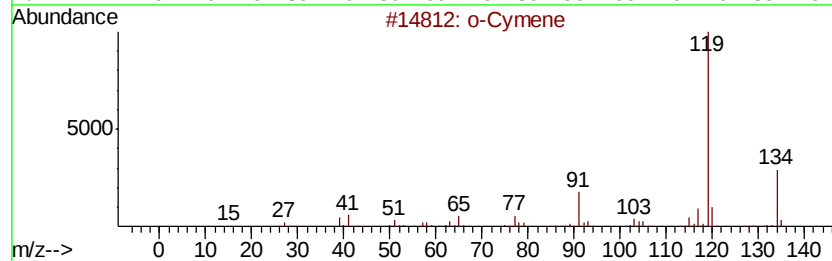
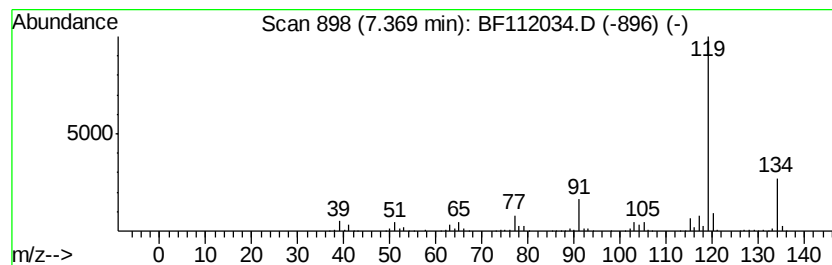
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 o-Cymene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	28.18 ng	1074890	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	97
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	97
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		p-Cymene	134	C10H14	000099-87-6	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
Acq On : 11 Jan 2019 23:49
Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CROSSWICKS-DRUM-1

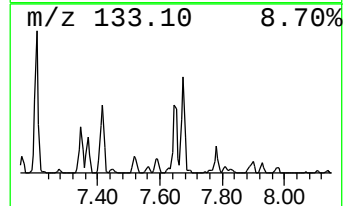
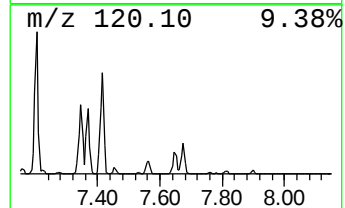
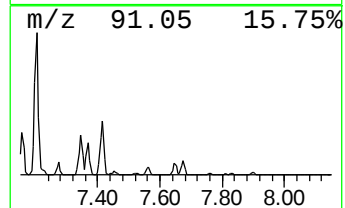
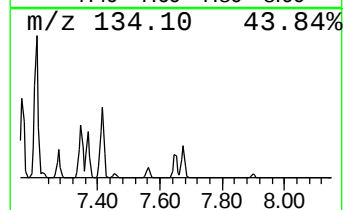
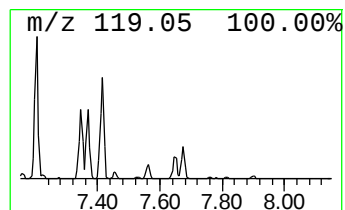
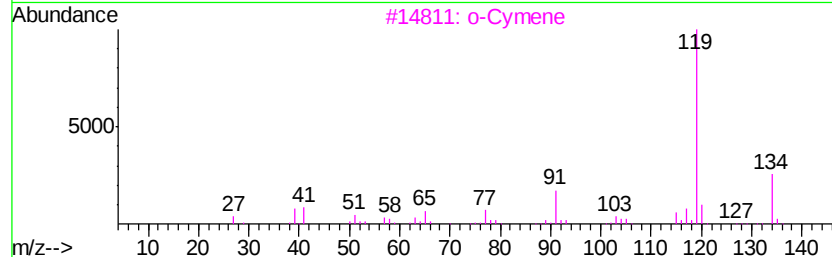
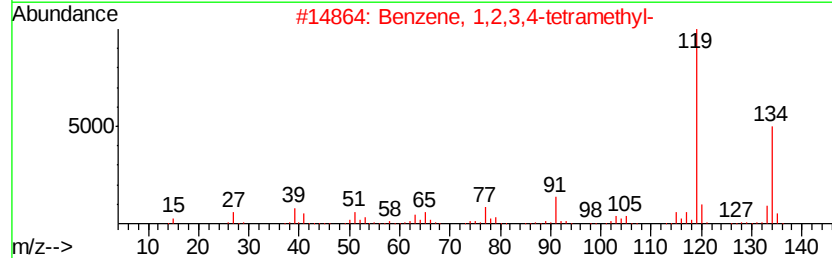
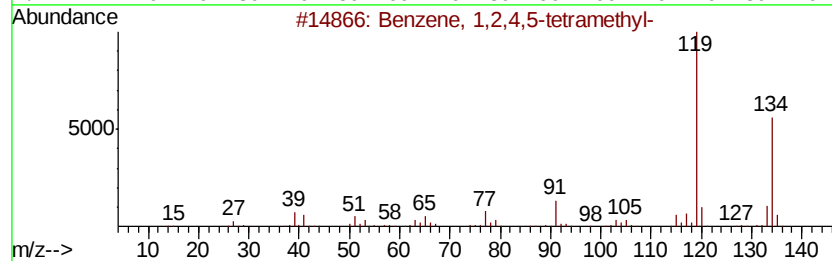
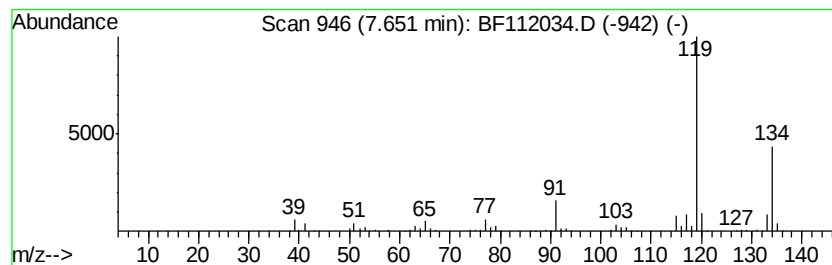
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	17.94 ng	487275	Napthalene-d8	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
Acq On : 11 Jan 2019 23:49
Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

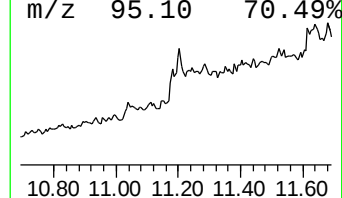
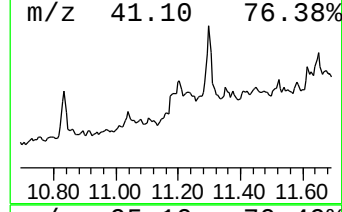
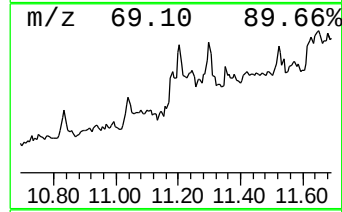
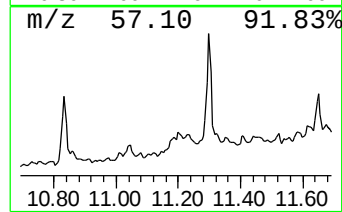
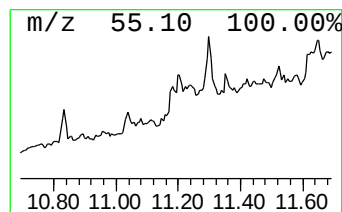
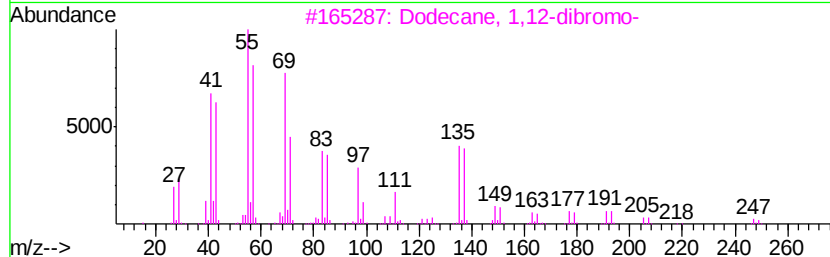
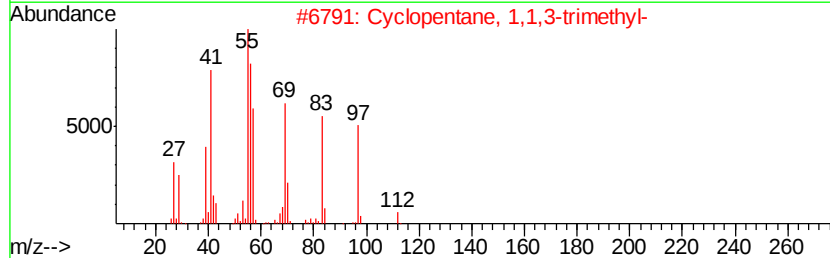
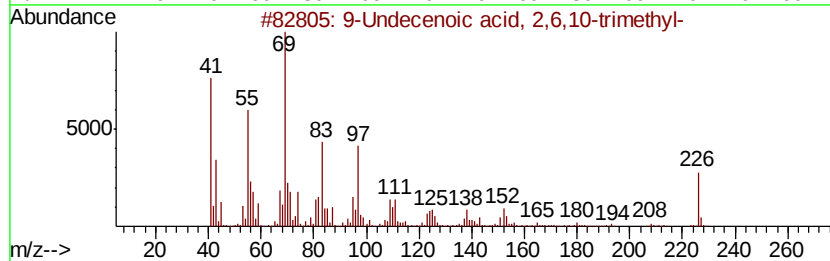
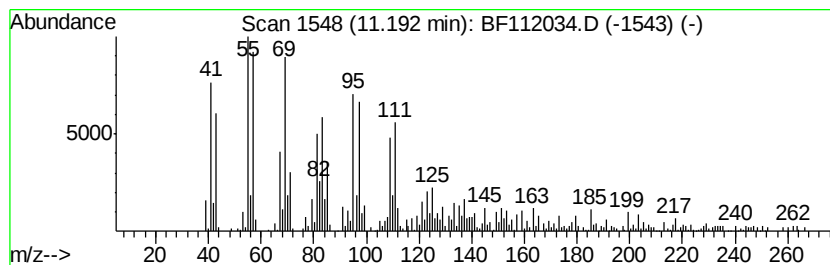
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 9-Undecenoic acid, 2,6,10-t... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.19	11.38 ng	508350	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Undecenoic acid, 2,6,10-trimet...	226	C14H26O2	1000131-86-2	52
2	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	45
3	Dodecane, 1,12-dibromo-	326	C12H24Br2	003344-70-5	38
4	Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	38
5	Cyclopentane, (2-methylpropyl)-	126	C9H18	003788-32-7	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
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Operator : JU/SJ
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ClientSampleId :
CROSSWICKS-DRUM-1

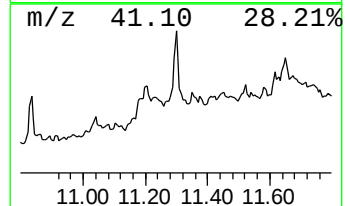
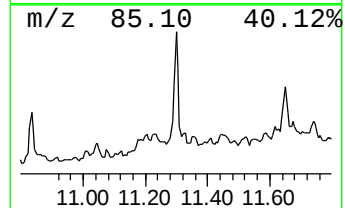
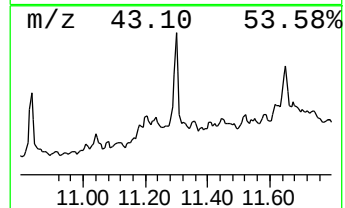
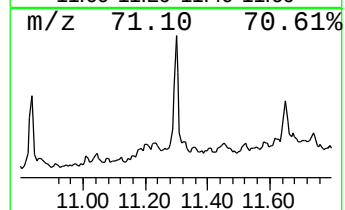
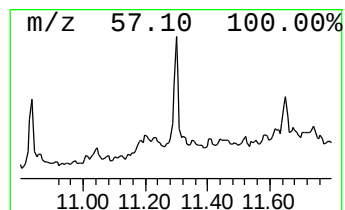
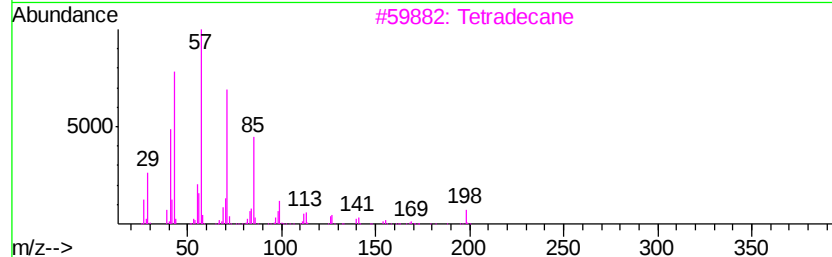
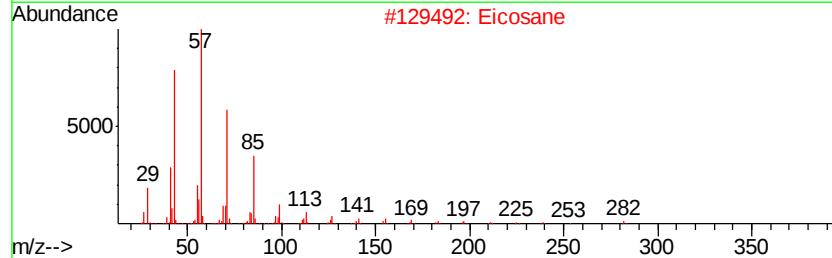
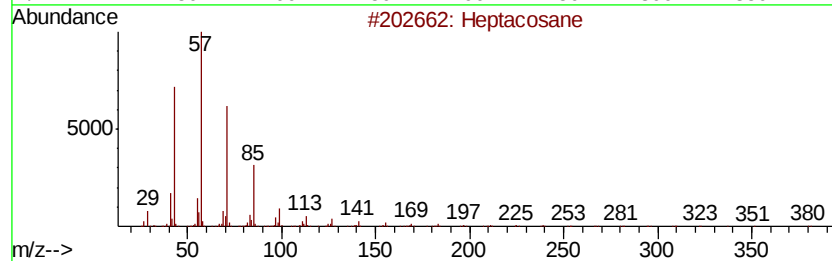
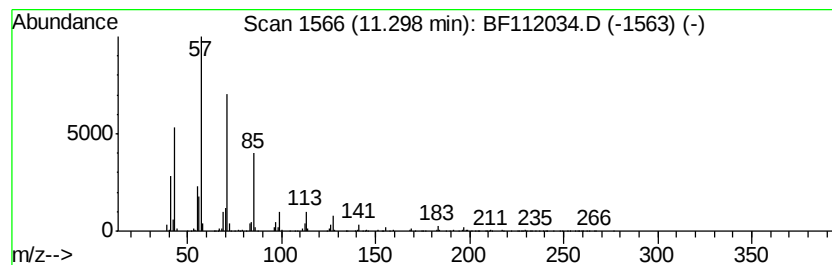
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heptacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	20.00 ng	893047	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	90
2		Eicosane	282	C20H42	000112-95-8	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Heptadecane	240	C17H36	000629-78-7	87
5		Tridecane, 4-methyl-	198	C14H30	026730-12-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
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Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

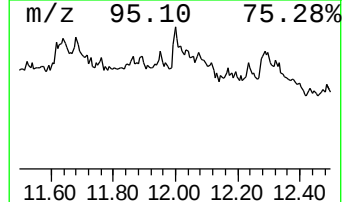
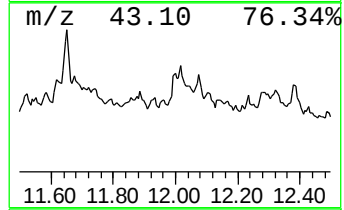
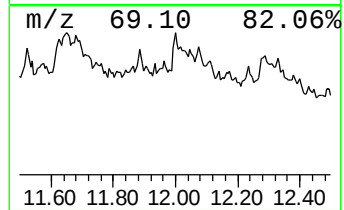
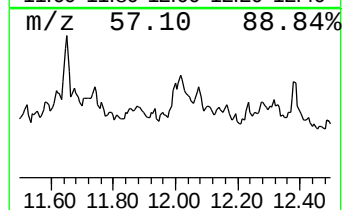
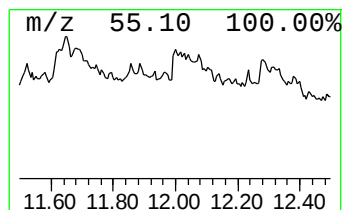
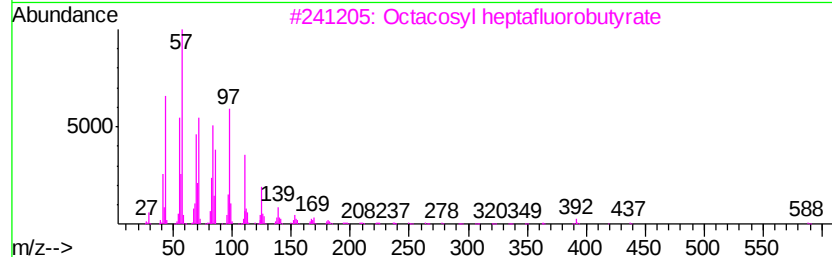
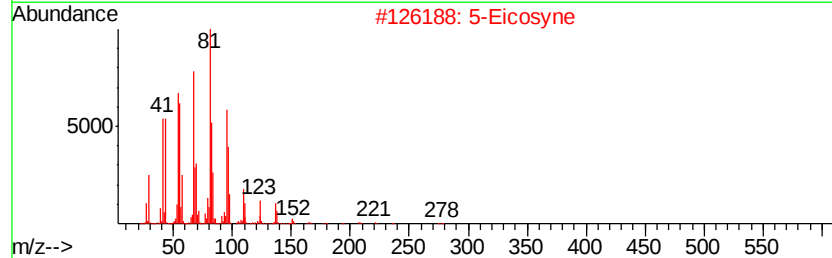
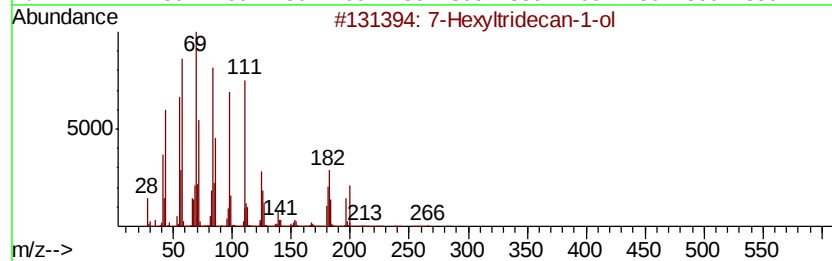
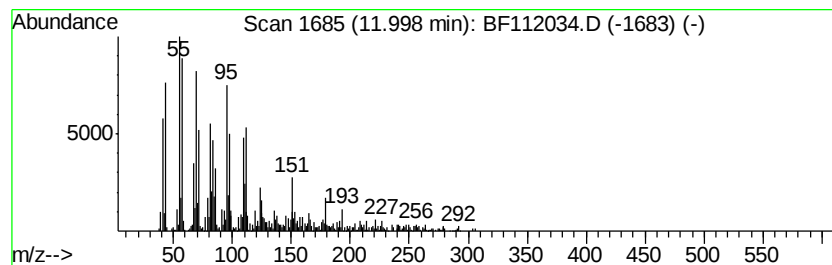
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ClientSampleId :
CROSSWICKS-DRUM-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 7-Hexyltridecan-1-ol Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.00	11.33 ng	505940	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Hexyltridecan-1-ol	284	C19H40O	1000197-13-2	51
2	5-Eicosyne	278	C20H38	074685-31-7	38
3	Octacosyl heptafluorobutyrate	606	C32H57F7O2	1000351-83-6	35
4	Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	35
5	Triacontyl heptafluorobutyrate	634	C34H61F7O2	1000351-83-2	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
Data File : BF112034.D
Acq On : 11 Jan 2019 23:49
Operator : JU/SJ
Sample : K1094-01 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

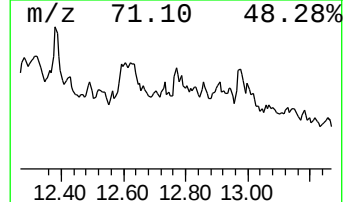
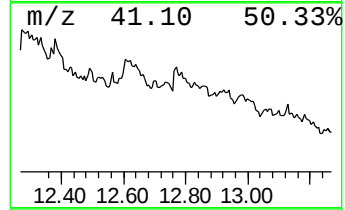
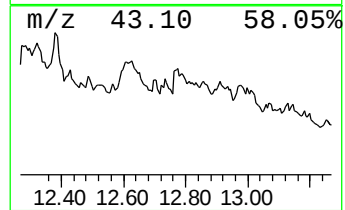
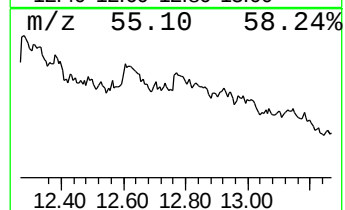
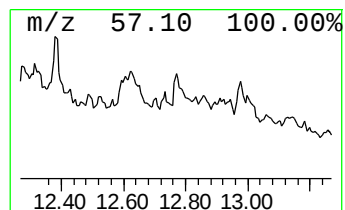
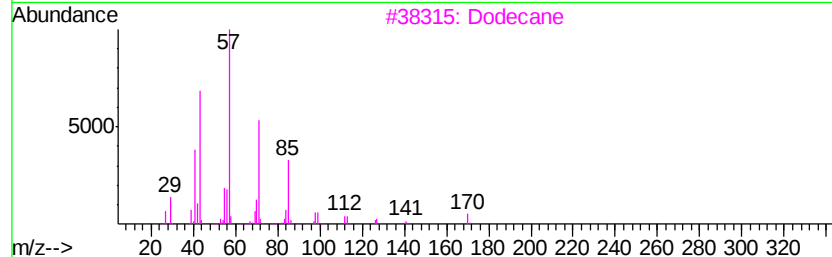
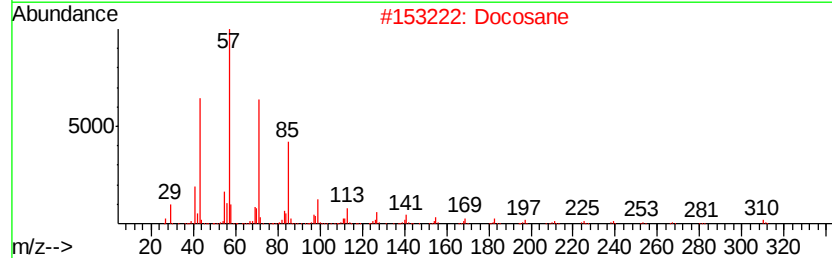
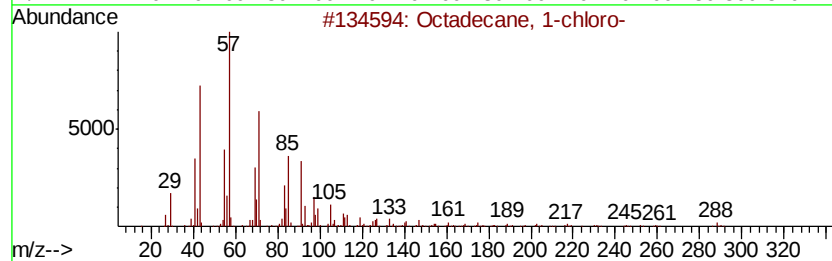
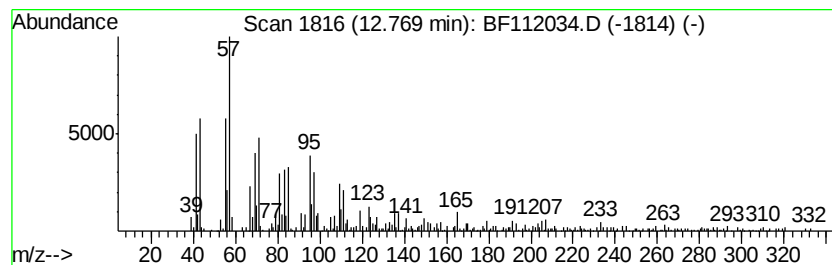
Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Octadecane, 1-chloro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	20.00 ng	678123	Chrysene-d12	14.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octadecane, 1-chloro-	288 C18H37Cl	003386-33-2 83
2		Docosane	310 C22H46	000629-97-0 74
3		Dodecane	170 C12H26	000112-40-3 70
4		2,6,6,10-Tetramethyl-undeca-8,10...	236 C15H24O2	098419-16-0 45
5		2-Dodecen-1-yl(-)succinic anhydride	266 C16H26O3	019780-11-1 45



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF011119\
 Data File : BF112034.D
 Acq On : 11 Jan 2019 23:49
 Operator : JU/SJ
 Sample : K1094-01 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CROSSWICKS-DRUM-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	6.42	202.3	ng	7716030	1	6.88	762742	20.0
Benzene, 1-ethyl-...	6.45	102.1	ng	3894980	1	6.88	762742	20.0
Indane	7.06	49.9	ng	1903760	1	6.88	762742	20.0
Benzene, 1,2-diet...	7.13	29.6	ng	1128220	1	6.88	762742	20.0
Benzene, 1-ethyl-...	7.20	102.2	ng	3895740	1	6.88	762742	20.0
Benzene, 1-methyl...	7.27	19.3	ng	736553	1	6.88	762742	20.0
Benzene, 2-ethyl-...	7.35	37.9	ng	1446430	1	6.88	762742	20.0
o-Cymene	7.37	28.2	ng	1074890	1	6.88	762742	20.0
Benzene, 1,2,4,5-...	7.65	17.9	ng	487275	2	8.16	543308	20.0
9-Undecenoic acid...	11.19	11.4	ng	508350	4	11.40	893047	20.0
Heptacosane	11.30	20.0	ng	893047	4	11.40	893047	20.0
7-Hexyltridecan-1-ol	12.00	11.3	ng	505940	4	11.40	893047	20.0
Octadecane, 1-chl...	12.77	20.0	ng	678123	5	14.04	678123	20.0