

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022415\
 Data File : BF077420.D
 Acq On : 24 Feb 2015 17:25
 Operator : IZ / TP
 Sample : PB81861BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB81861BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF021915.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	1.024	10	12	20	rVB6	12588	32382	1.45%	0.207%
2	1.595	58	62	65	rBV2	18252	43431	1.94%	0.277%
3	1.744	74	75	84	rVB	23817	41729	1.87%	0.267%
4	1.870	84	86	94	rBV	61147	82908	3.71%	0.530%
5	5.093	365	368	373	rBV	520056	664359	29.72%	4.244%
6	5.584	409	411	414	rBV	771674	1062008	47.51%	6.785%
7	6.853	519	522	524	rBV	1133136	1126361	50.39%	7.196%
8	7.013	533	536	538	rBV	807088	1125248	50.34%	7.189%
9	7.299	559	561	564	rBV	473729	418928	18.74%	2.676%
10	7.482	575	577	580	rVB	975248	1172332	52.45%	7.489%
11	7.985	619	621	624	rBV	751174	863016	38.61%	5.513%
12	8.876	697	699	702	rBV	639824	570612	25.53%	3.645%
13	10.225	814	817	819	rBV	2226082	1928687	86.29%	12.321%
14	10.751	860	863	866	rVB	23965	44284	1.98%	0.283%
15	11.048	887	889	892	rBV	518605	657417	29.41%	4.200%
16	12.031	973	975	978	rBV	901078	845851	37.84%	5.404%
17	12.899	1048	1051	1053	rBV	675577	708558	31.70%	4.527%
18	14.579	1196	1198	1201	rVB	35355	34643	1.55%	0.221%
19	14.900	1223	1226	1229	rBV	2291789	2235125	100.00%	14.279%
20	16.248	1342	1344	1347	rBV	562425	780547	34.92%	4.986%
21	16.945	1394	1405	1408	rBV7	8096	46504	2.08%	0.297%
22	18.168	1509	1512	1515	rBV	537011	802517	35.90%	5.127%
23	20.214	1685	1691	1702	rBV	75137	365788	16.37%	2.337%

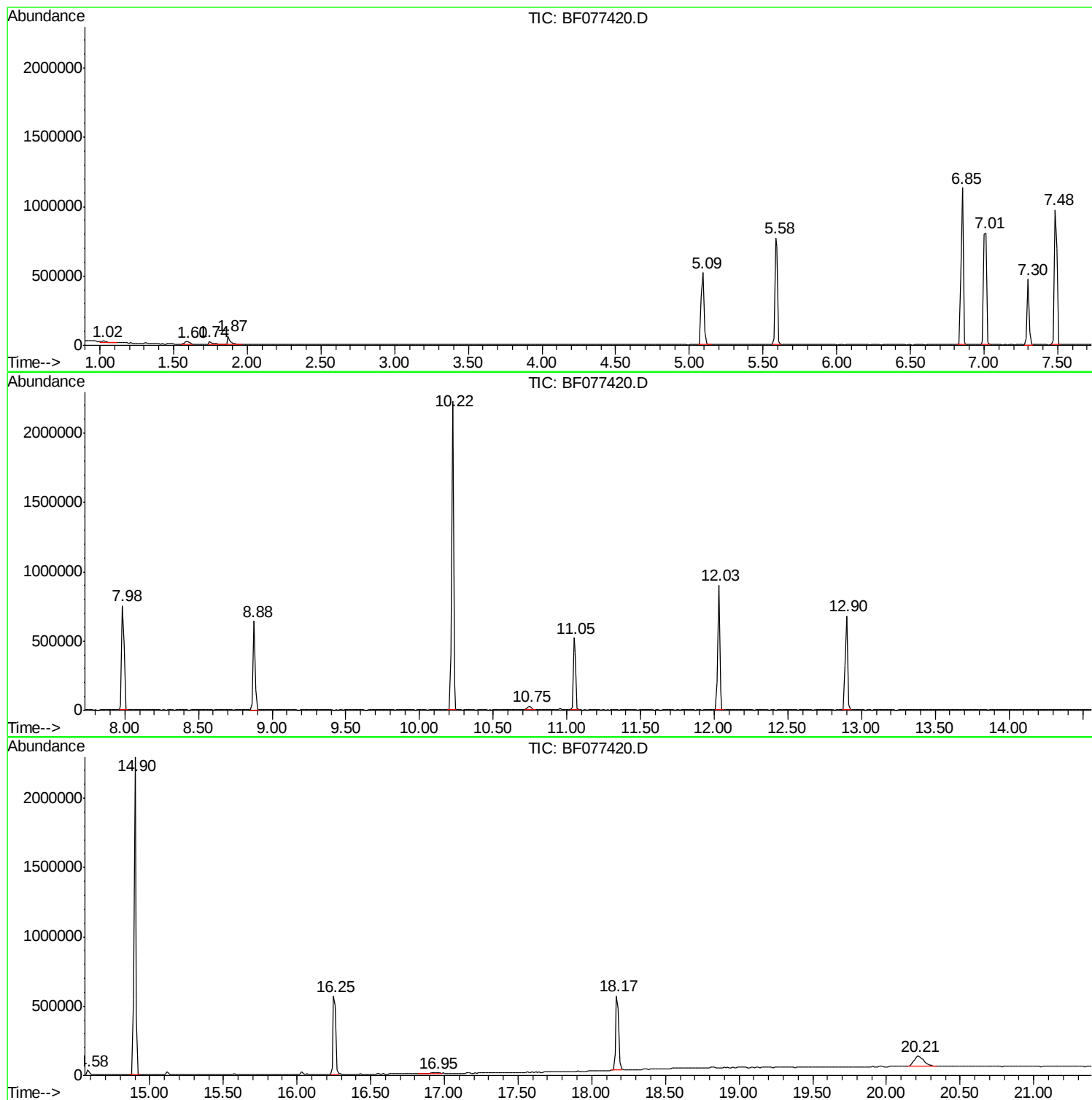
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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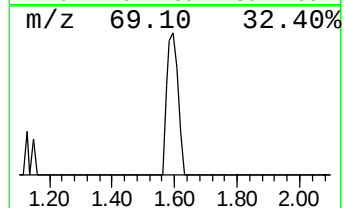
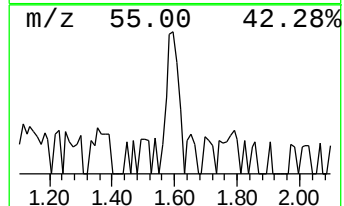
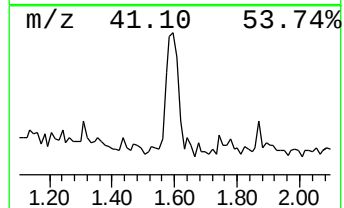
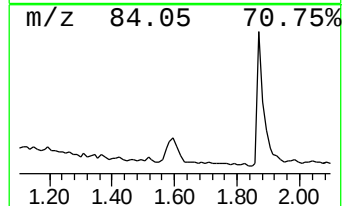
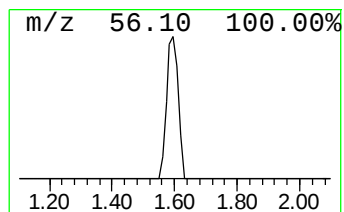
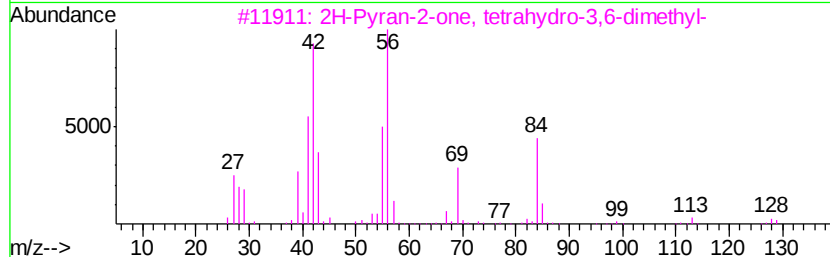
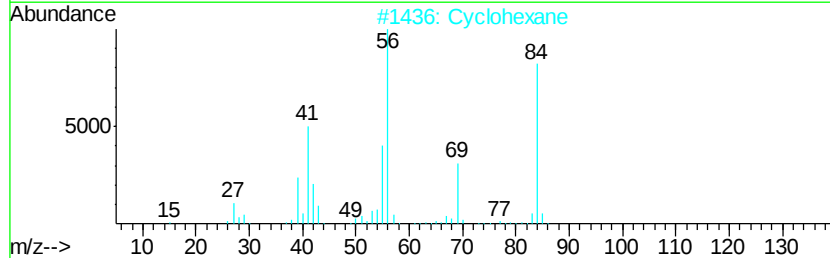
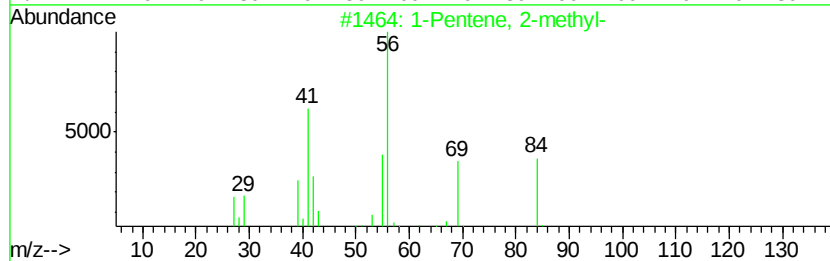
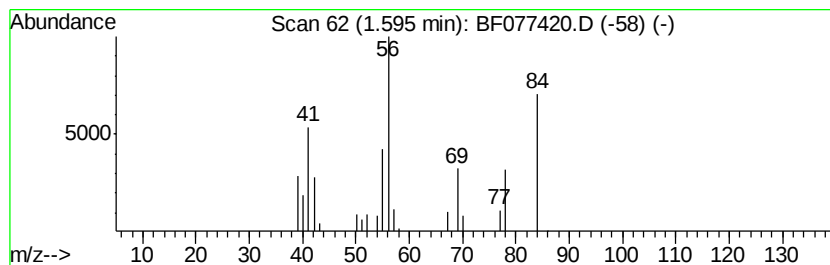
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.60	2.07 ng	43431	1,4-Dichlorobenzene-d4	7.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
2			Cyclohexane	84	C6H12	000110-82-7	58
3			2H-Pyran-2-one, tetrahydro-3,6-d...	128	C7H12O2	003720-22-7	50
4			Oxirane, (1-methylbutyl)-	114	C7H14O	053229-39-3	47
5			3-Buten-1-ol, 2-methyl-	86	C5H10O	004516-90-9	38



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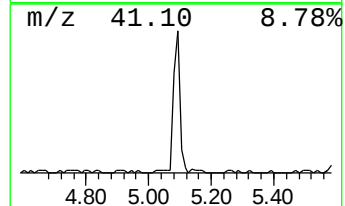
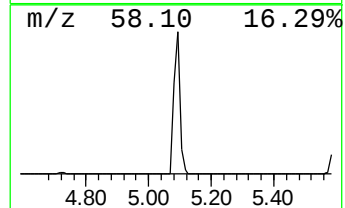
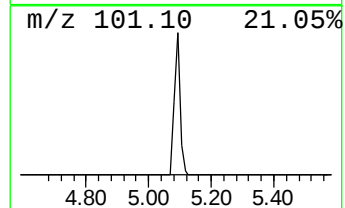
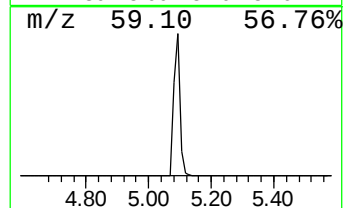
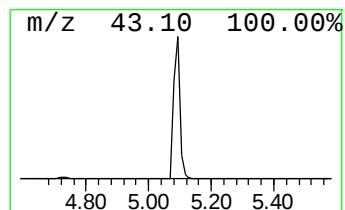
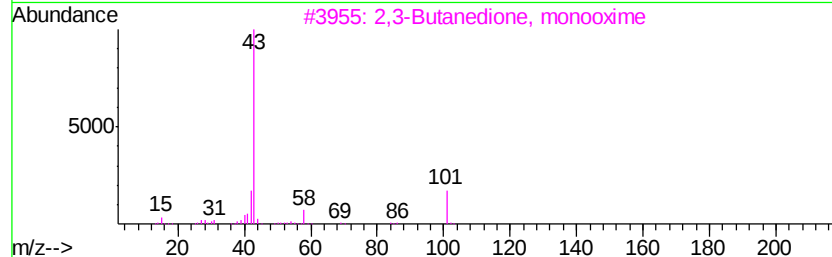
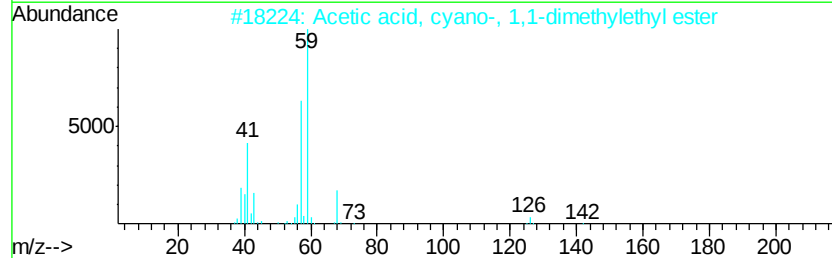
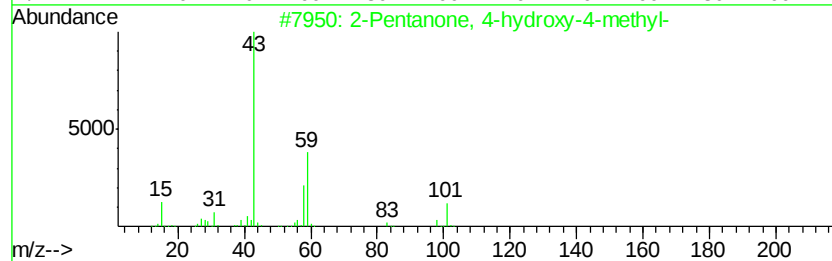
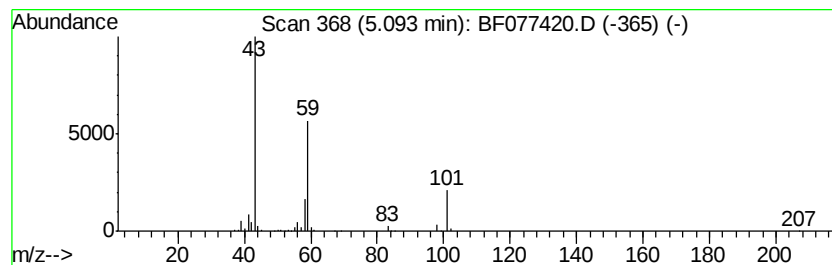
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	31.72 ng	664359	1,4-Dichlorobenzene-d4	7.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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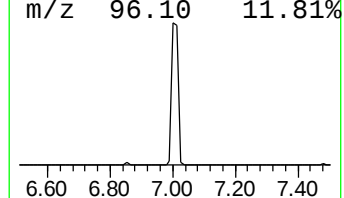
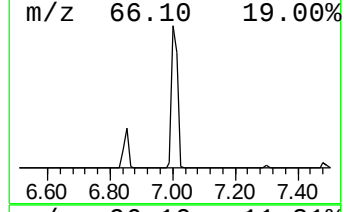
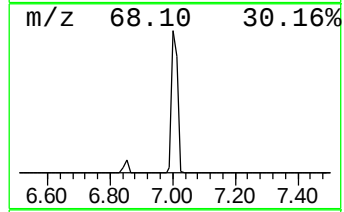
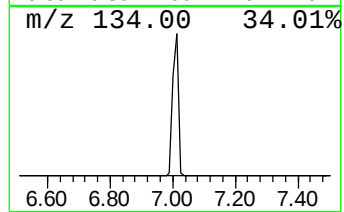
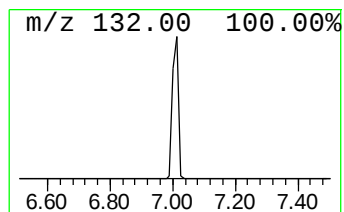
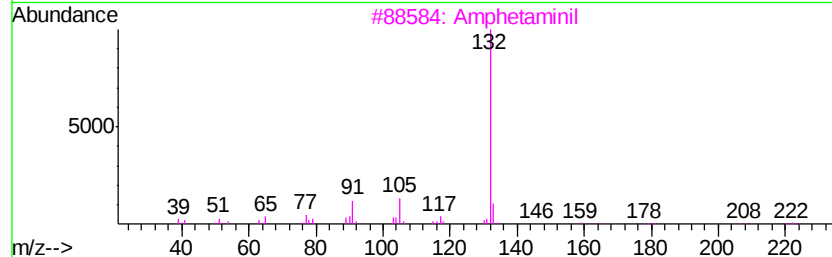
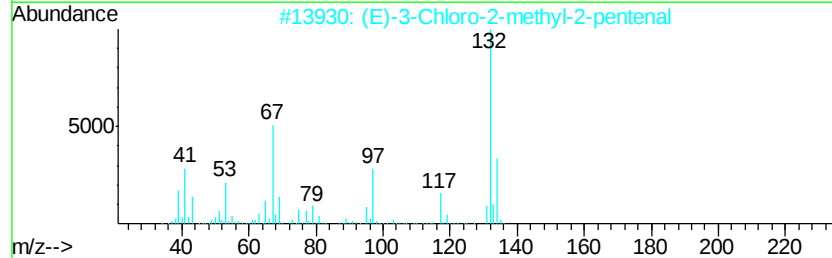
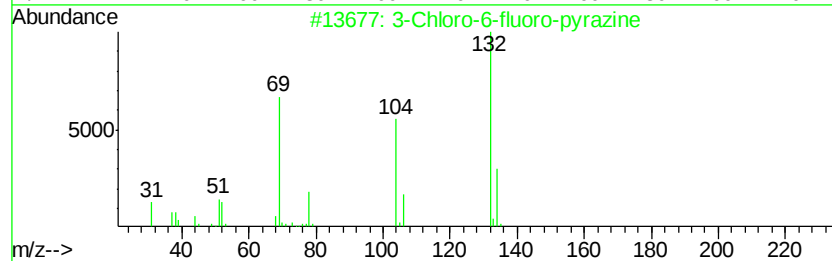
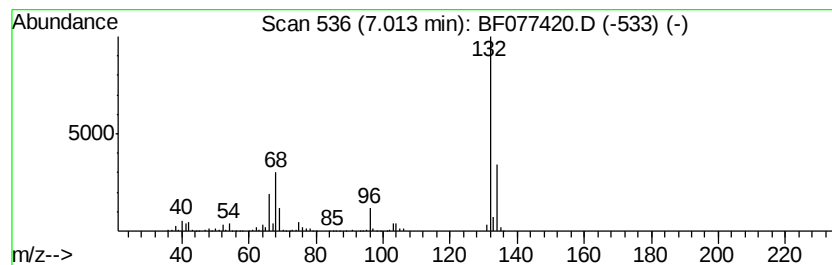
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Peak Number 4 unknown7.01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.01	53.72 ng	1125250	1,4-Dichlorobenzene-d4	7.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	40
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	33
3		Amphetaminil	250	C17H18N2	017590-01-1	25
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		Xenon	132	Xe	007440-63-3	9



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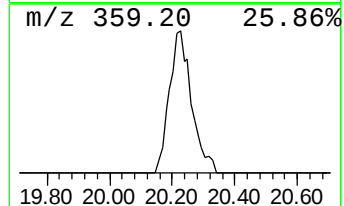
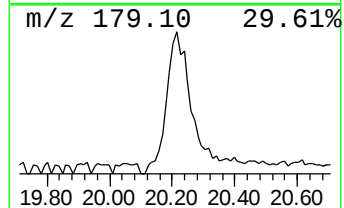
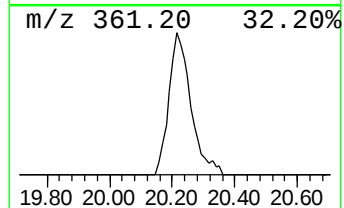
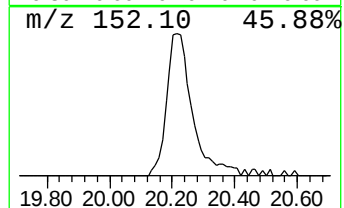
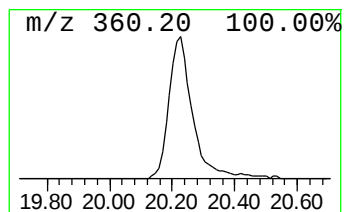
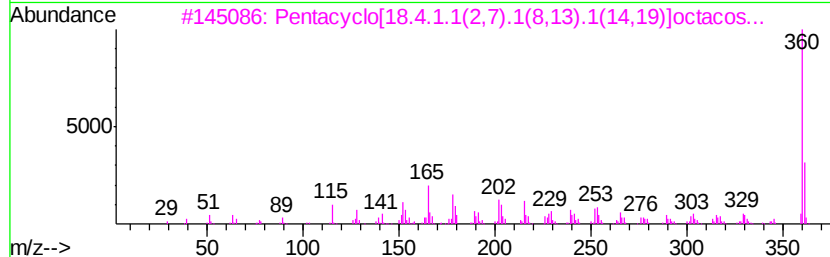
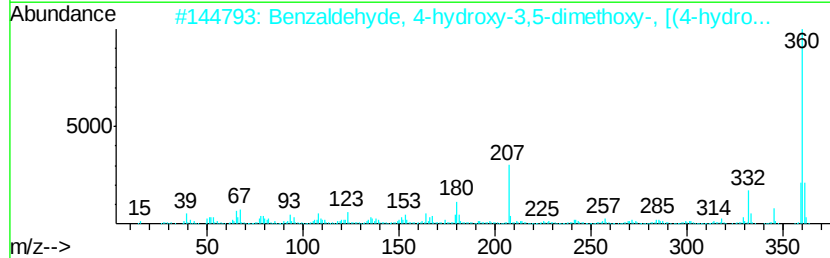
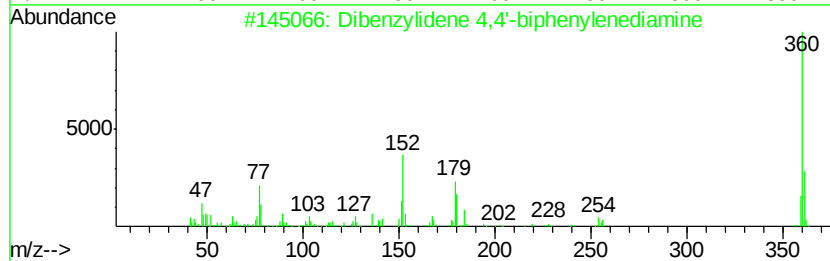
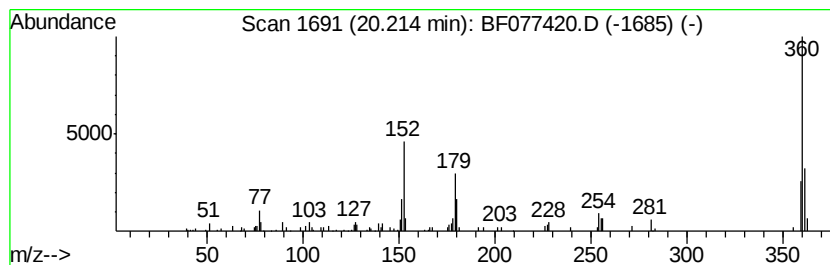
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Peak Number 6 Dibenzylidene 4,4'-biphenyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.21	9.12 ng	365788	Perylene-d12	18.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzylidene 4,4'-biphenylenedi...	360	C26H20N2	006311-48-4	93
2		Benzaldehyde, 4-hydroxy-3,5-dime...	360	C18H20N2O6	014414-32-5	47
3		Pentacyclo[18.4.1.1(2,7).1(8,13)...]	360	C28H24	101077-60-5	47
4		N,N'-di-2-Naphthyl-p-phenylenedi...	360	C26H20N2	000093-46-9	46
5		1,3,2-Dioxaborolane, bis(spiro[4...	360	C20H13B06	1000150-23-7	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	1.60	2.1	ng	43431	1	7.30	418928	20.0
2-Pentanone, 4-hy...	5.09	31.7	ng	664359	1	7.30	418928	20.0
unknown7.01	7.01	53.7	ng	1125250	1	7.30	418928	20.0
Dibenzylidene 4,4...	20.21	9.1	ng	365788	6	18.17	802517	20.0