

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
 Data File : BF080174.D
 Acq On : 25 Jun 2015 19:09
 Operator : TP/IZ
 Sample : G2778-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NNESIDEWALL-062415

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:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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	5.510	209	212	214	rBV	254935	367018	25.45%	3.508%
2	5.899	244	246	249	rBV	925035	943285	65.42%	9.016%
3	6.299	279	281	284	rBV	25511	29778	2.07%	0.285%
4	6.893	331	333	336	rBV	838890	972404	67.44%	9.294%
5	7.064	345	348	350	rBV	1042869	1052754	73.01%	10.062%
6	7.293	366	368	371	rVB	219258	190137	13.19%	1.817%
7	7.453	380	382	385	rBV	959189	769577	53.37%	7.356%
8	7.853	414	417	419	rBV	790484	656983	45.56%	6.280%
9	8.585	479	481	484	rBV	324235	261994	18.17%	2.504%
10	9.659	573	575	578	rVB	1220078	1100432	76.31%	10.518%
11	10.048	607	609	612	rBV	363300	299680	20.78%	2.864%
12	10.356	634	636	639	rVB	352841	301298	20.89%	2.880%
13	10.768	669	672	674	rBV	13179	16912	1.17%	0.162%
14	11.156	703	706	708	rBV2	584375	805019	55.83%	7.695%
15	11.248	712	714	719	rVB	19740	21655	1.50%	0.207%
16	11.865	765	768	770	rBV	344965	331047	22.96%	3.164%
17	12.071	783	786	790	rVB2	8392	15456	1.07%	0.148%
18	12.139	790	792	794	rBV	13382	17480	1.21%	0.167%
19	13.534	912	914	917	rBV	1558215	1441964	100.00%	13.783%
20	14.585	1003	1006	1015	rBV	51923	105606	7.32%	1.009%
21	14.791	1022	1024	1027	rBV	377918	374033	25.94%	3.575%
22	15.762	1107	1109	1114	rBV2	22232	33200	2.30%	0.317%
23	16.654	1184	1187	1190	rBV	225824	354518	24.59%	3.389%

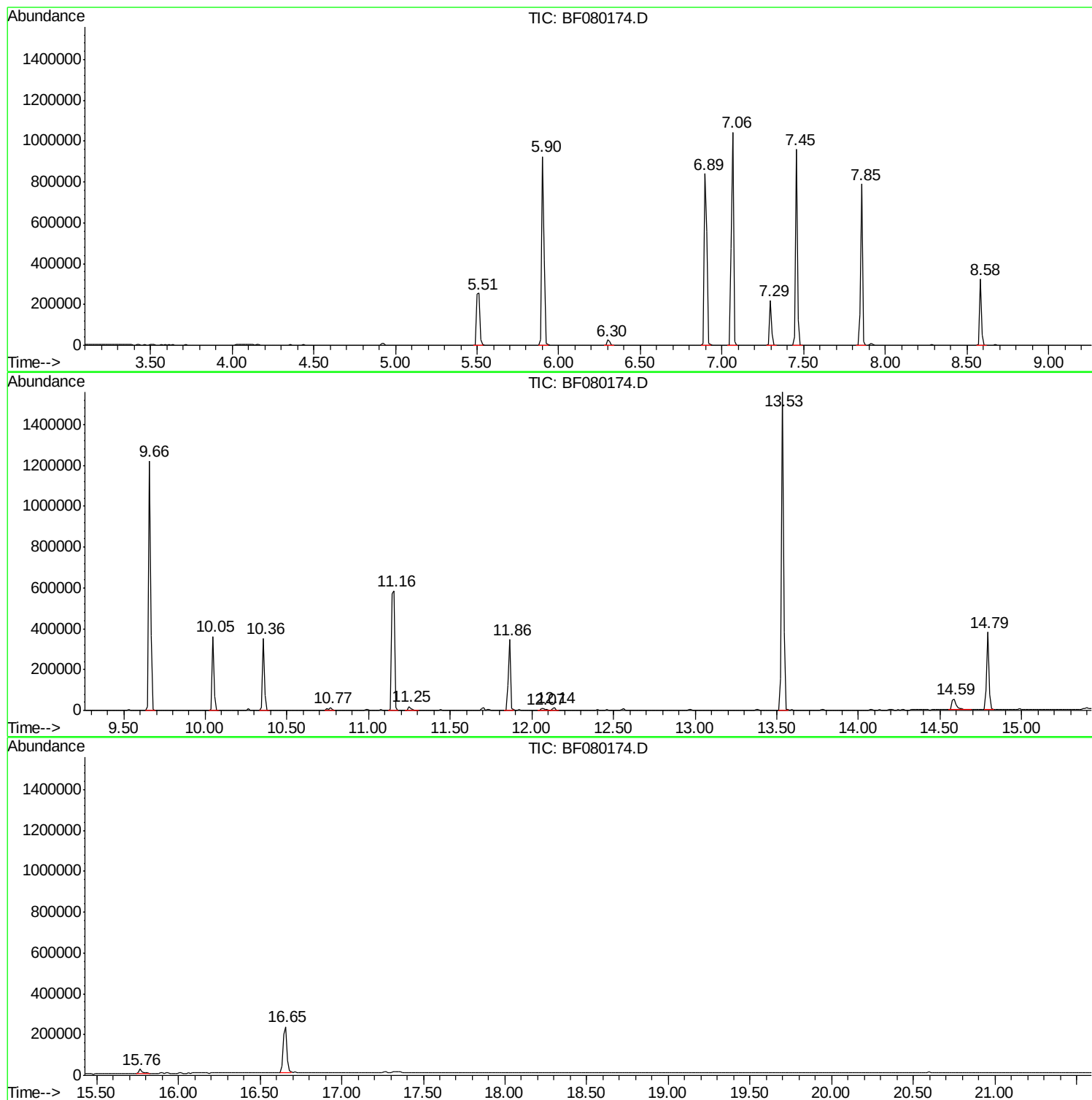
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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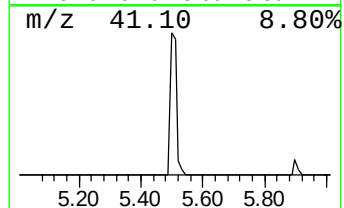
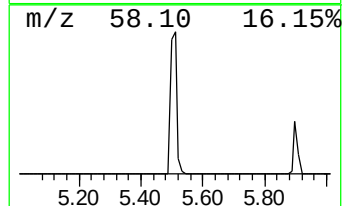
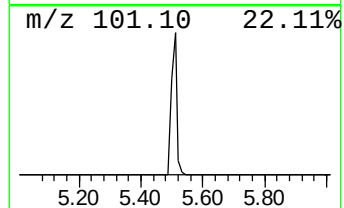
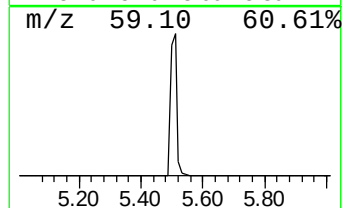
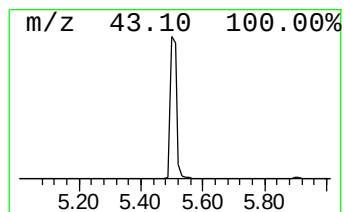
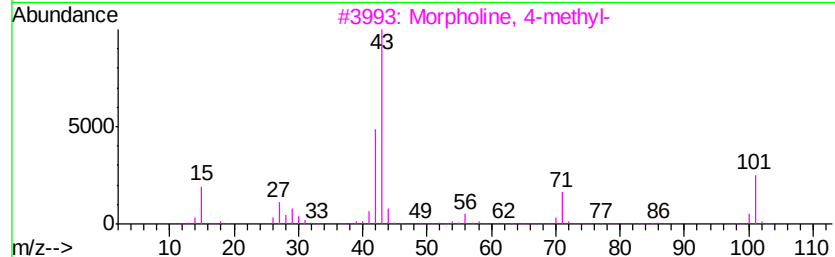
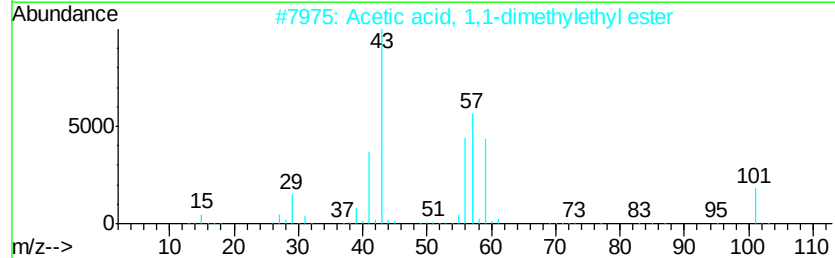
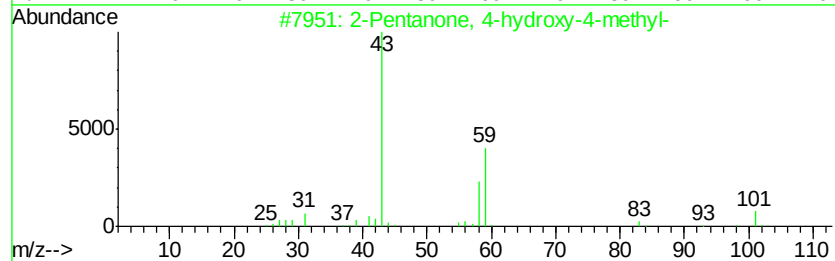
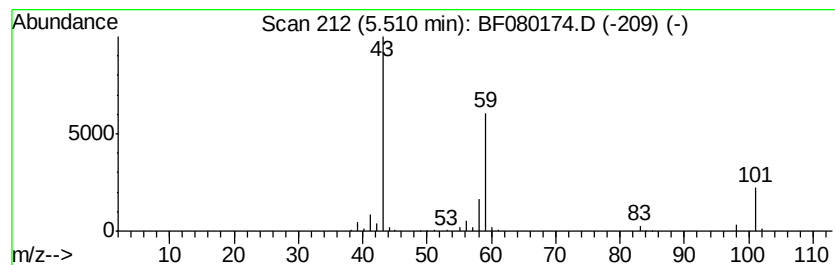
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	38.61 ng	367018	1,4-Dichlorobenzene-d4	7.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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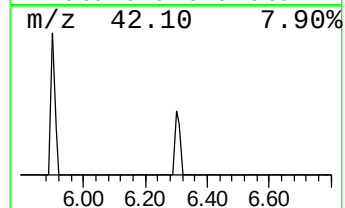
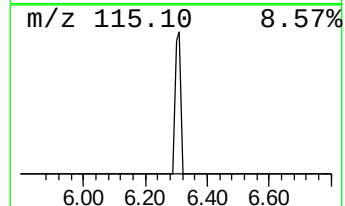
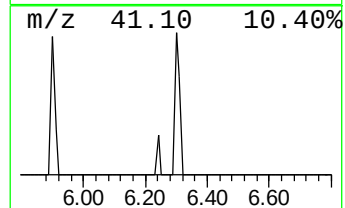
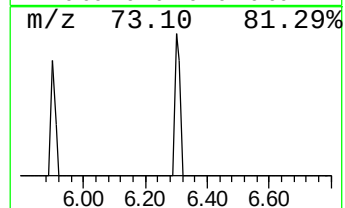
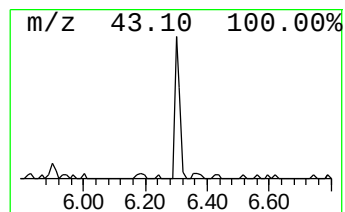
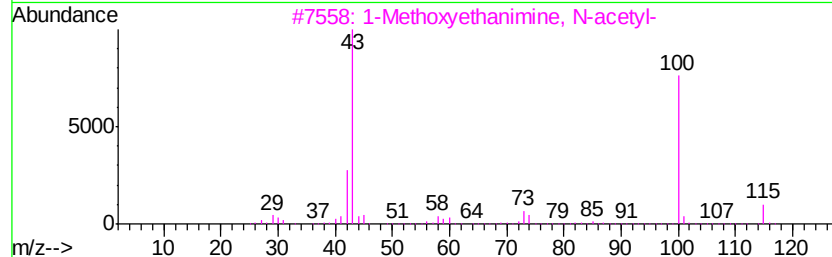
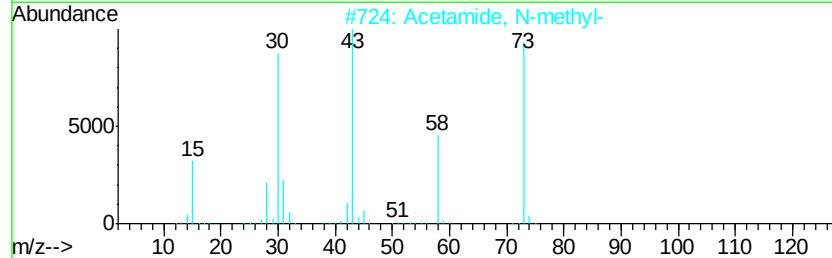
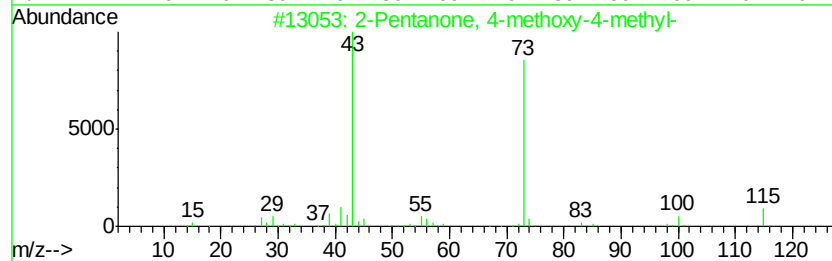
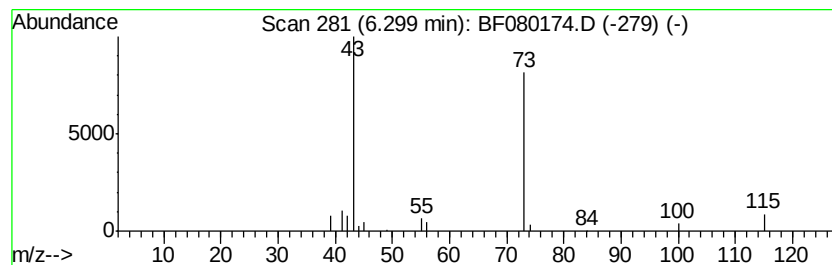
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Acetamide, N-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.30	3.13 ng	29778	1,4-Dichlorobenzene-d4	7.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	50
3		1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9



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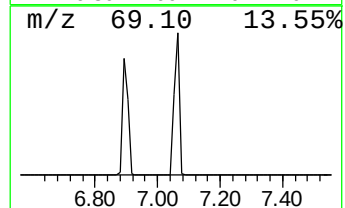
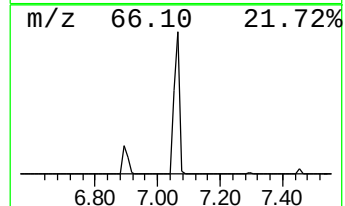
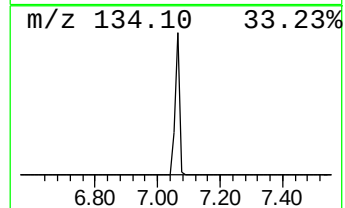
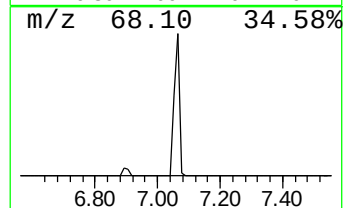
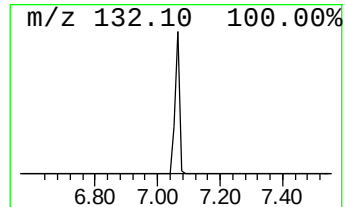
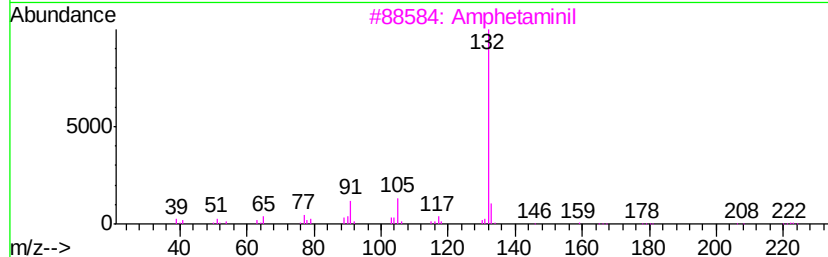
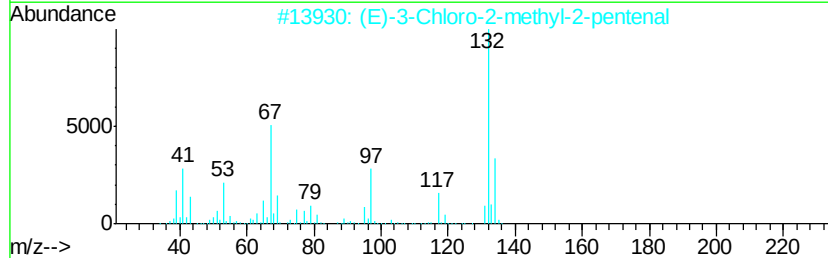
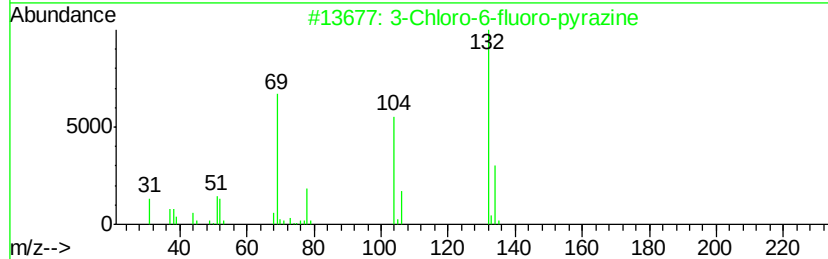
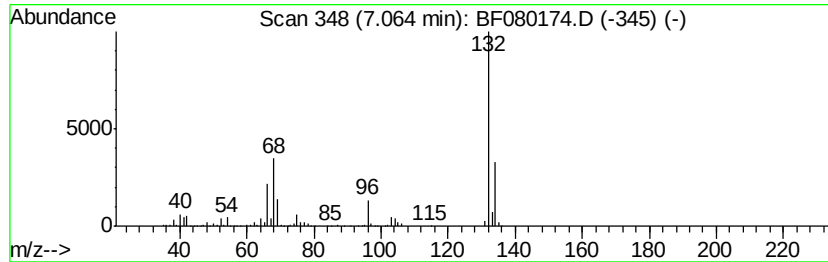
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Peak Number 3 unknown7.06 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	110.74 ng	1052750	1,4-Dichlorobenzene-d4	7.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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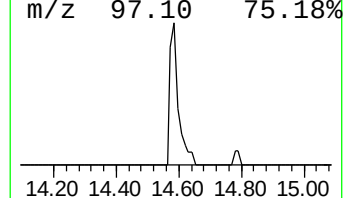
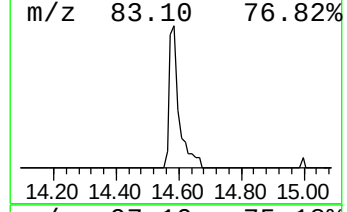
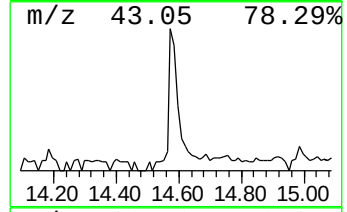
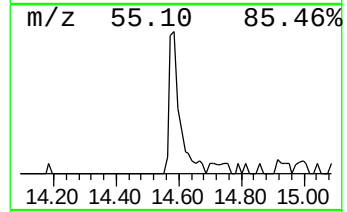
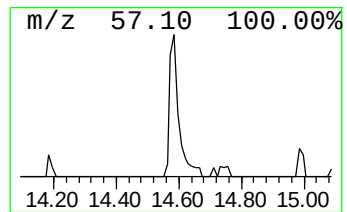
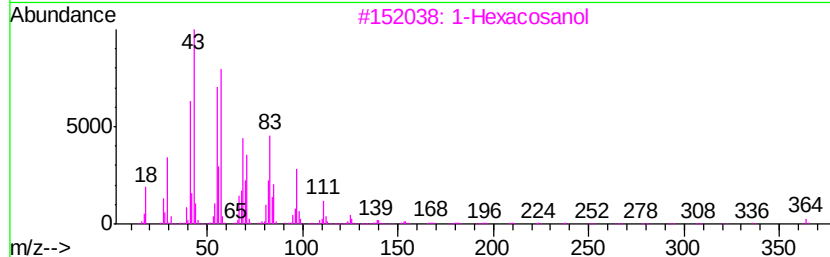
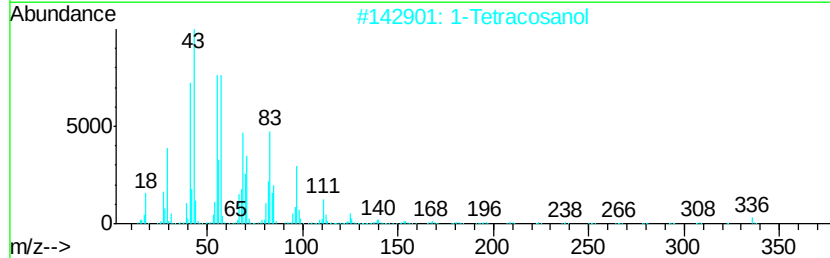
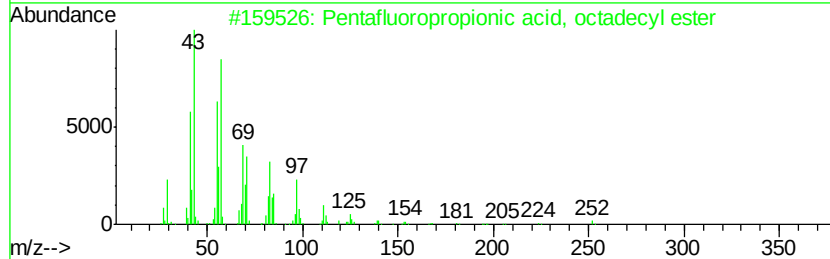
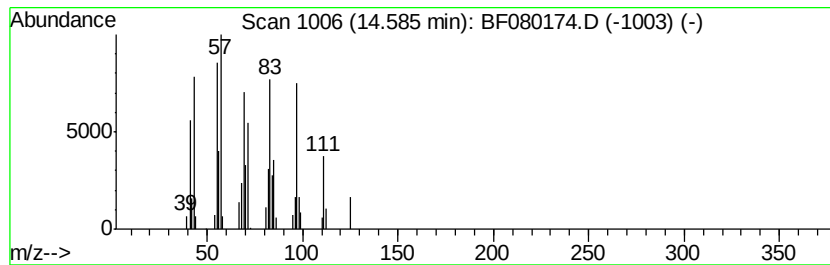
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Peak Number 5 Pentafluoropropionic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.59	5.65 ng	105606	Chrysene-d12		14.79	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentafluoropropionic acid, octadec...	416	C21H37F5O2	1000280-07-7	87
2		1-Tetracosanol	354	C24H50O	000506-51-4	64
3		1-Hexacosanol	382	C26H54O	000506-52-5	64
4		Acetic acid, trifluoro-, tetradec...	310	C16H29F3O2	006222-02-2	50
5		1H-Imidazole, 4,5-dihydro-2-(1-m...	112	C6H12N2	040029-86-5	46



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.51	38.6	ng	367018	1	7.29	190137	20.0
Acetamide, N-methyl-	6.30	3.1	ng	29778	1	7.29	190137	20.0
unknown7.06	7.06	110.7	ng	1052750	1	7.29	190137	20.0
Pentafluoropropio...	14.59	5.7	ng	105606	5	14.79	374033	20.0