

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
 Data File : BF077391.D
 Acq On : 20 Feb 2015 19:40
 Operator : TP/IZ
 Sample : PB81820BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB81820BL

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-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	1.515	52	55	57	rVB	6486932	7597549	100.00%	24.992%
2	1.630	63	65	75	rVB2	41534	129241	1.70%	0.425%
3	1.767	75	77	86	rVV	179372	403033	5.30%	1.326%
4	1.892	86	88	110	rVB	689369	1222382	16.09%	4.021%
5	5.127	369	371	379	rBV	377779	435901	5.74%	1.434%
6	5.619	412	414	417	rBV	1585742	1752805	23.07%	5.766%
7	6.876	521	524	527	rBV	1897835	1736115	22.85%	5.711%
8	7.036	536	538	541	rBV	2018240	1949094	25.65%	6.412%
9	7.333	562	564	566	rBV	507841	491069	6.46%	1.615%
10	7.516	578	580	583	rBV	1781568	1567004	20.63%	5.155%
11	8.019	621	624	626	rBV	1364095	1177968	15.50%	3.875%
12	8.910	699	702	704	rBV	560354	660857	8.70%	2.174%
13	10.259	817	820	822	rBV	1967623	2659069	35.00%	8.747%
14	11.082	890	892	895	rBV	891794	842077	11.08%	2.770%
15	12.065	975	978	980	rBV	1511697	1619882	21.32%	5.329%
16	12.934	1051	1054	1056	rBV	707850	907651	11.95%	2.986%
17	14.922	1226	1228	1231	rBV	2350891	3367703	44.33%	11.078%
18	16.225	1340	1342	1345	rBV	727472	1009860	13.29%	3.322%
19	18.008	1496	1498	1501	rVB	671595	870696	11.46%	2.864%

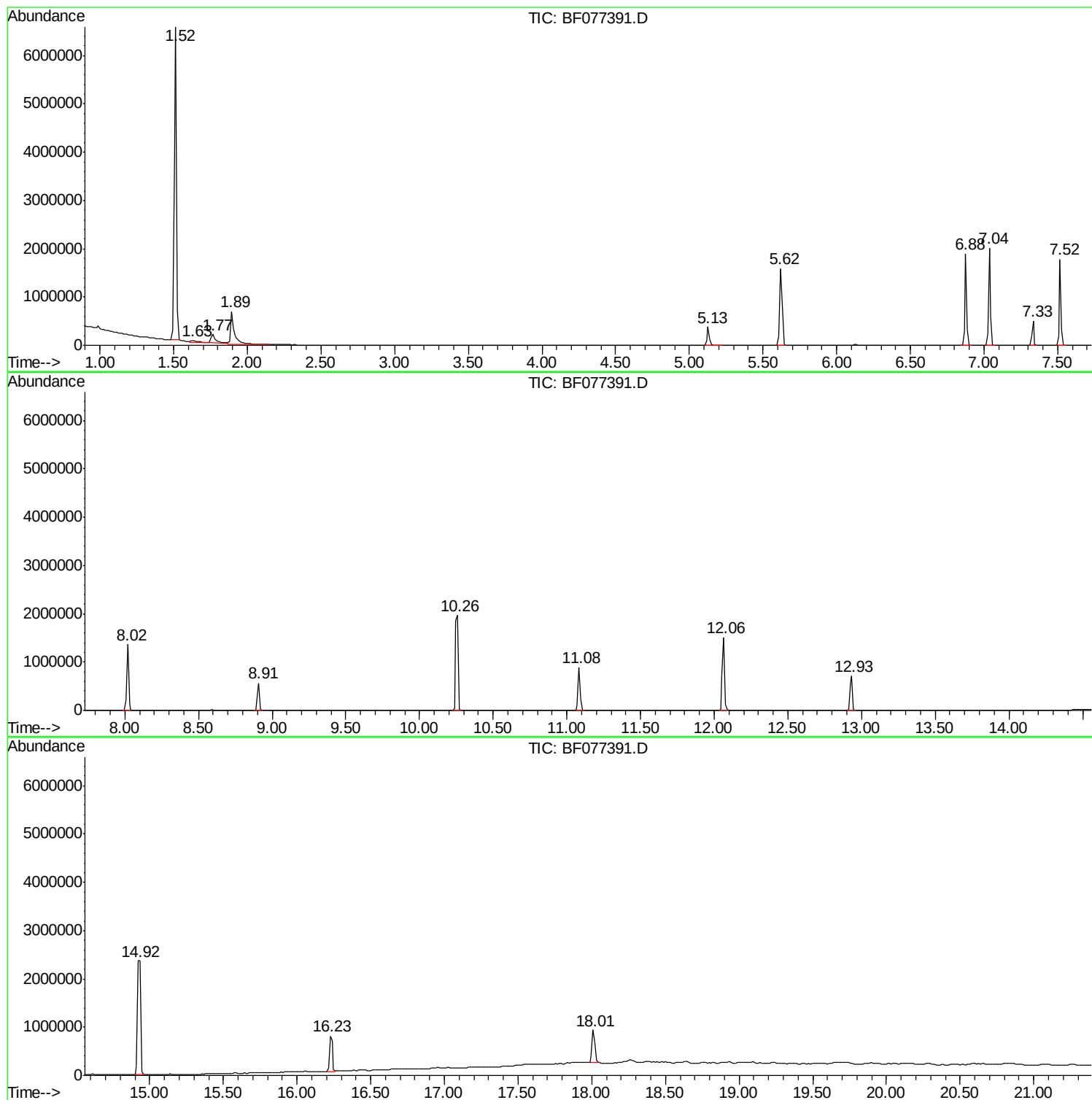
Sum of corrected areas: 30399956

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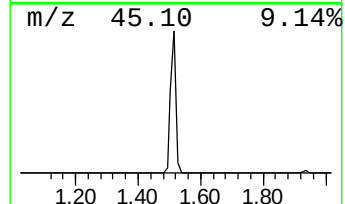
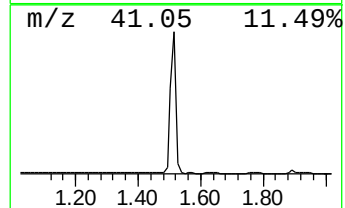
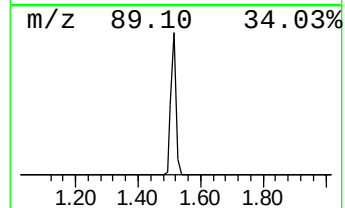
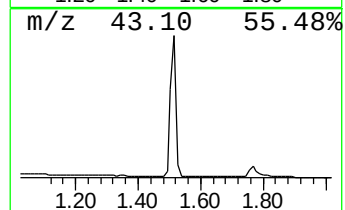
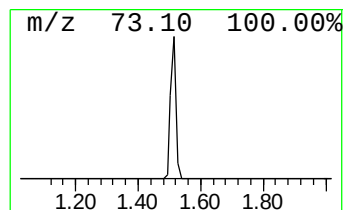
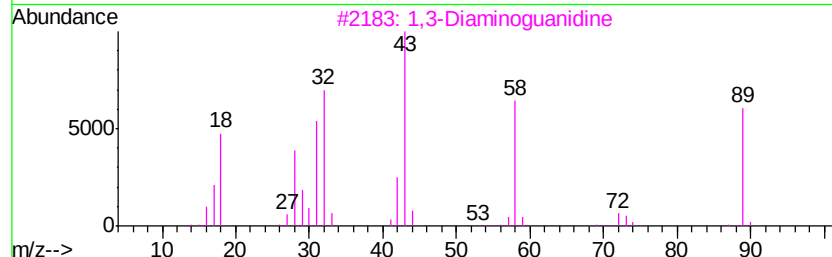
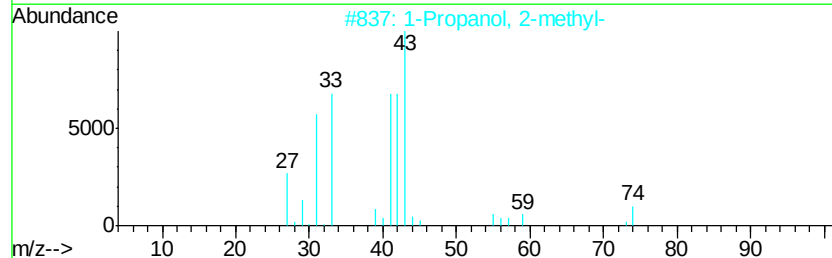
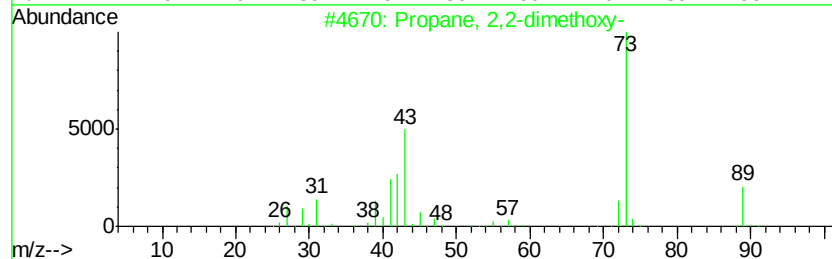
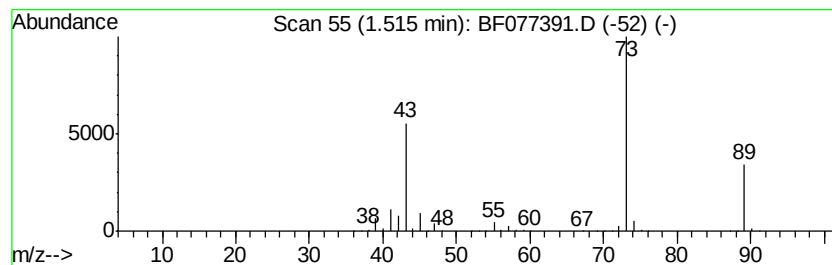
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	309.43 ng	7597550	1,4-Dichlorobenzene-d4	7.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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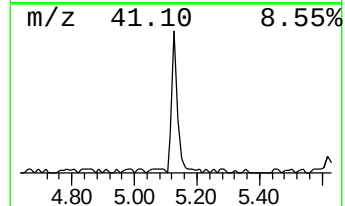
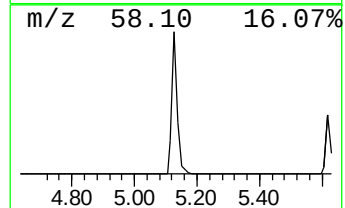
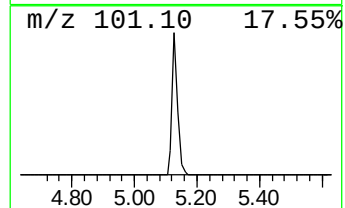
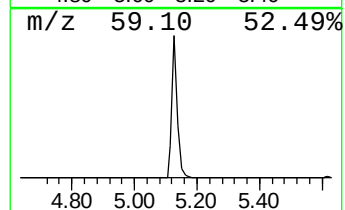
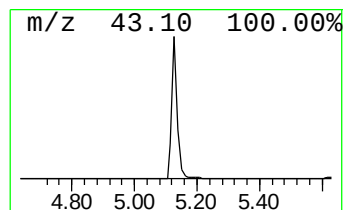
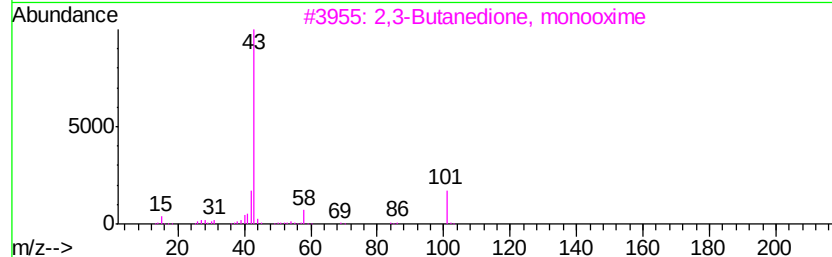
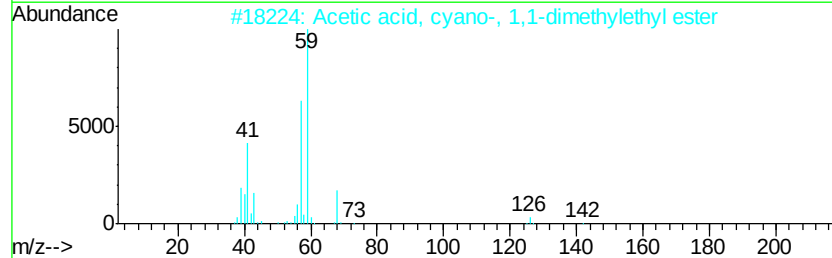
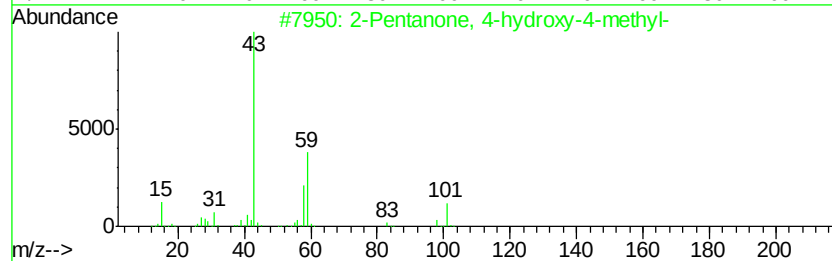
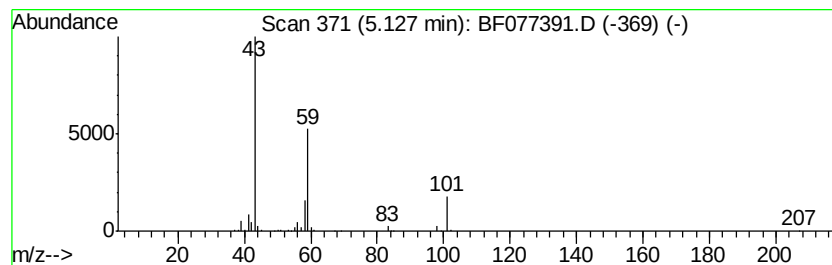
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Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	17.75 ng	435901	1,4-Dichlorobenzene-d4	7.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9



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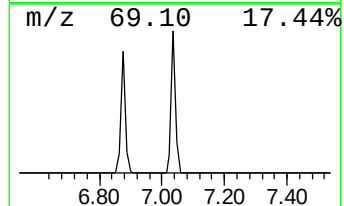
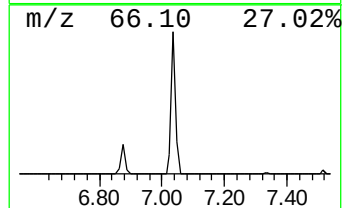
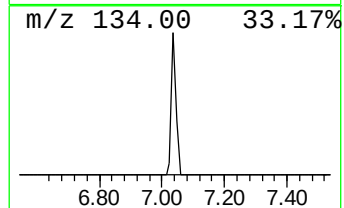
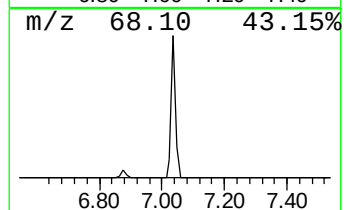
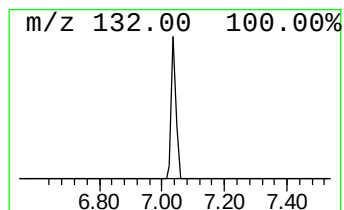
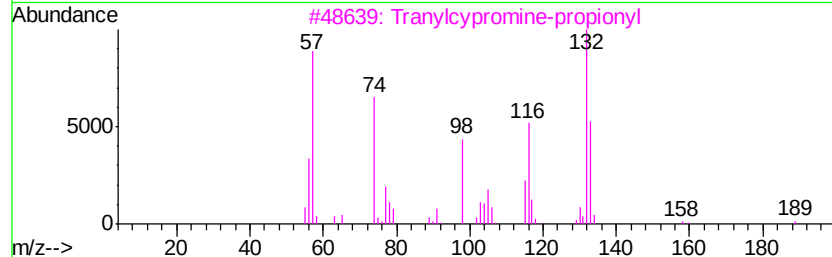
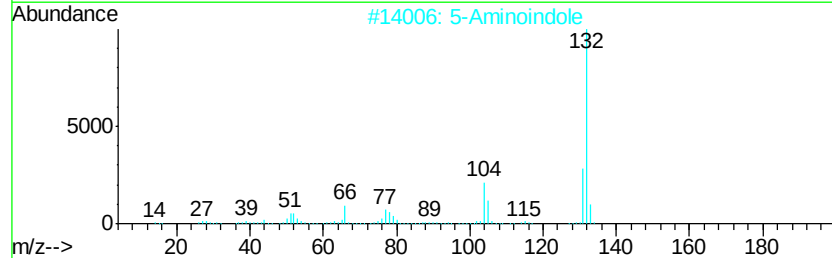
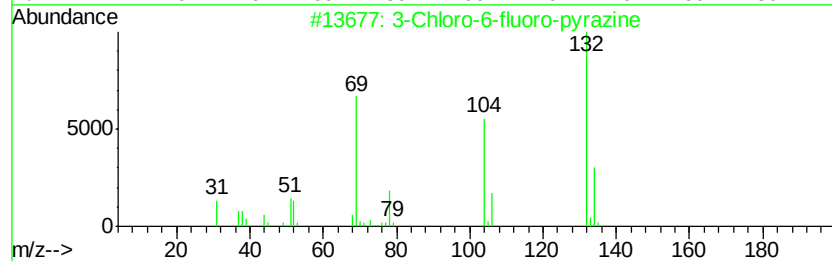
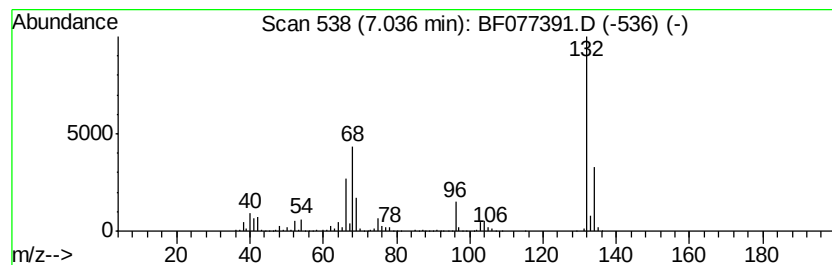
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Peak Number 6 unknown7.04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	79.38 ng	1949090	1,4-Dichlorobenzene-d4	7.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...			132	C10H12	028749-81-7	9
5	2H-Cyclopenta[d]pyridazine, 2-me...			132	C8H8N2	022291-85-6	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-dime...	1.52	309.4	ng	7597550	1	7.33	491069	20.0
2-Pentanone, 4-hy...	5.13	17.8	ng	435901	1	7.33	491069	20.0
unknown7.04	7.04	79.4	ng	1949090	1	7.33	491069	20.0