

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072415\
 Data File : BF080644.D
 Acq On : 25 Jul 2015 00:14
 Operator : TP/UM
 Sample : G3079-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12-7-22-15

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-BF072415.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	4.430	204	205	209	rVB2	12444	19331	1.30%	0.181%
2	5.081	261	262	265	rBV	189459	220709	14.83%	2.070%
3	5.504	297	299	302	rBV	799922	685813	46.07%	6.432%
4	5.927	335	336	338	rBV	45374	43539	2.92%	0.408%
5	6.533	387	389	392	rBV	356481	401550	26.97%	3.766%
6	6.693	400	403	405	rBV	1217569	1182244	79.41%	11.087%
7	6.933	421	424	426	rBV	340689	324386	21.79%	3.042%
8	7.082	435	437	440	rBV	908171	981027	65.90%	9.200%
9	7.493	470	473	475	rBV	891912	890063	59.79%	8.347%
10	8.213	534	536	539	rBV	376311	384741	25.84%	3.608%
11	9.287	628	630	633	rBV	1163389	1464704	98.39%	13.736%
12	9.973	687	690	692	rBV	544589	445045	29.89%	4.174%
13	10.762	756	759	761	rBV2	774792	936561	62.91%	8.783%
14	11.459	817	820	822	rBV	524884	434066	29.16%	4.071%
15	13.071	958	961	963	rBV	1083513	1488733	100.00%	13.962%
16	14.202	1058	1060	1063	rBV	427541	411437	27.64%	3.859%
17	15.825	1199	1202	1205	rBV	276774	349017	23.44%	3.273%

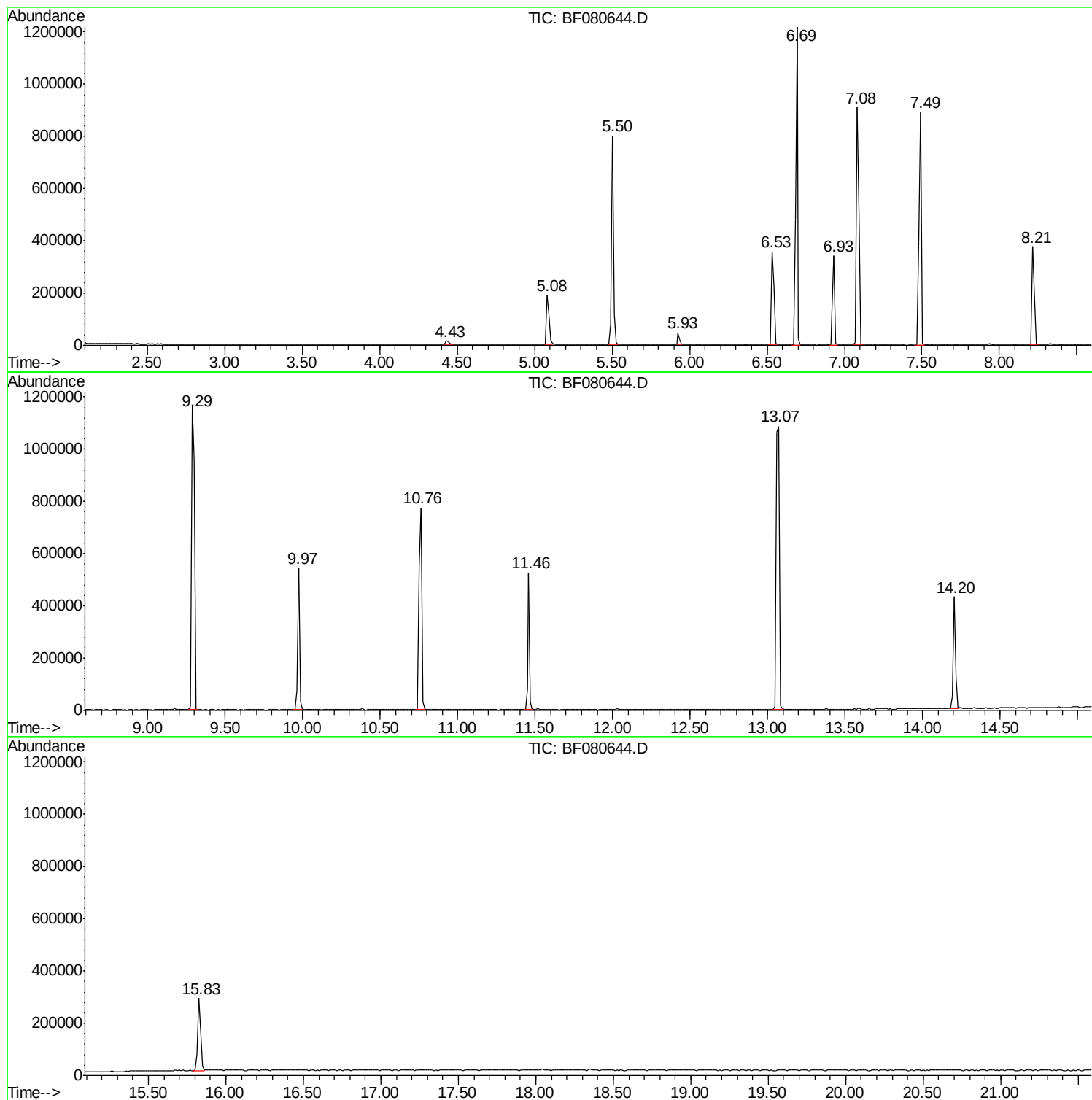
Sum of corrected areas: 10662966

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072415.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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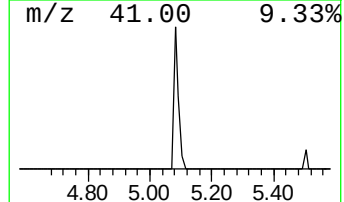
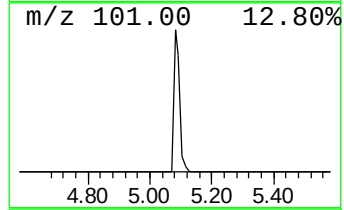
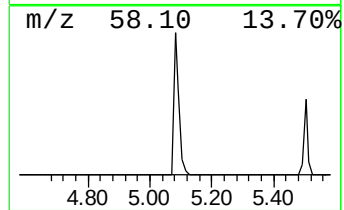
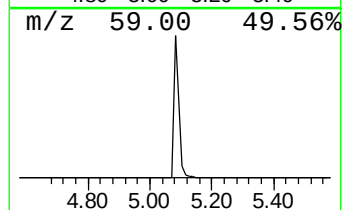
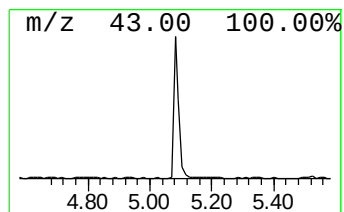
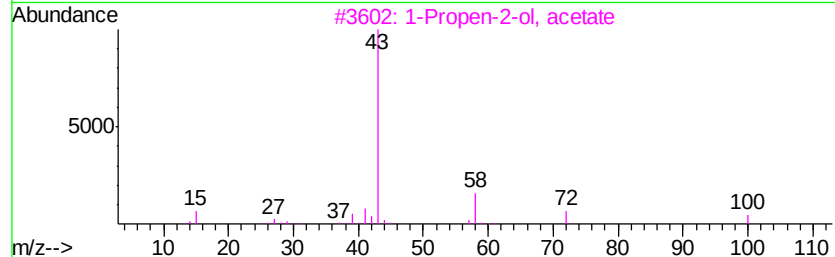
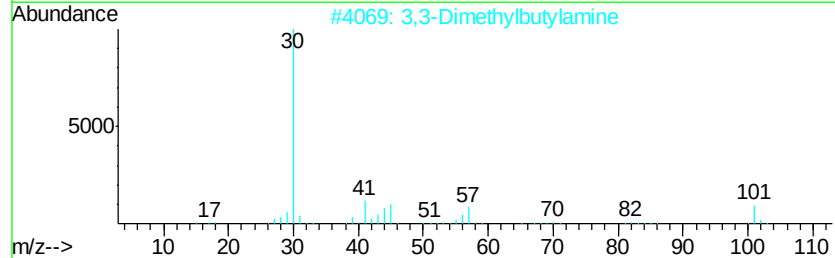
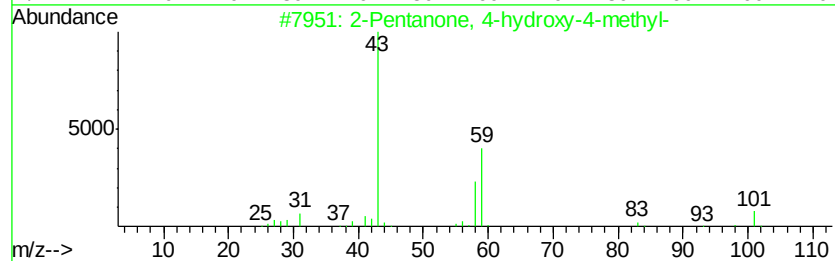
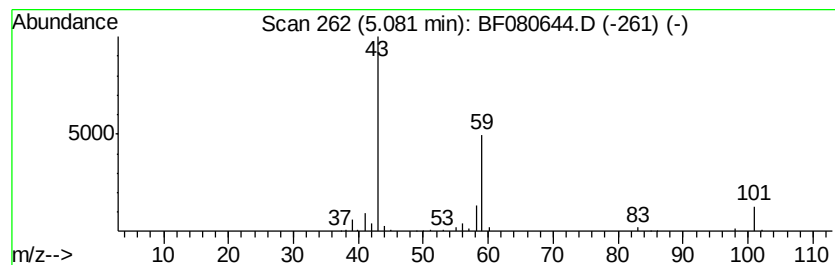
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.08	13.61 ng	220709	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3,3-Dimethylbutylamine	101	C6H15N	015673-00-4	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	9
5		2-Pentanone, 5-(acetyloxy)-	144	C7H12O3	005185-97-7	9



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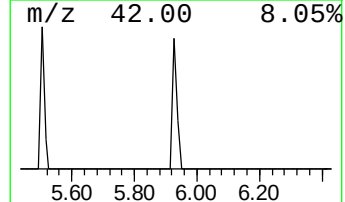
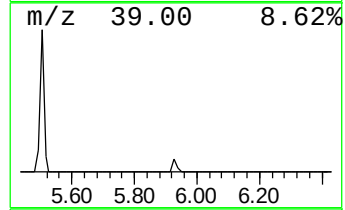
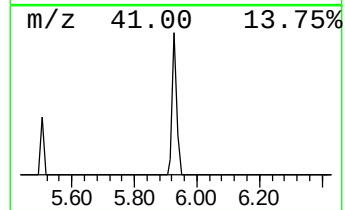
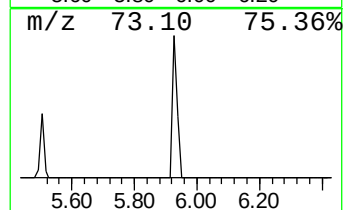
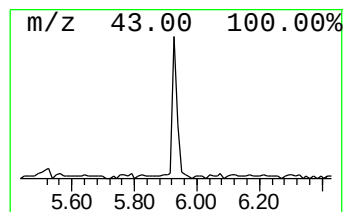
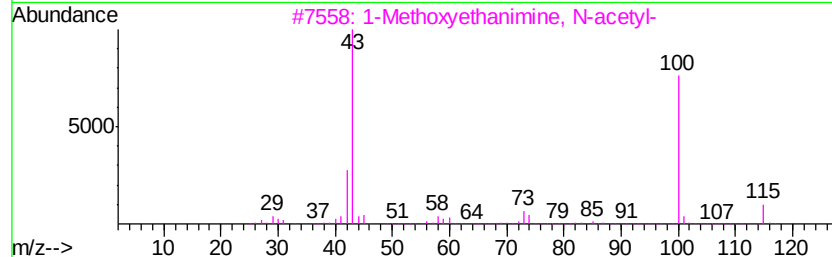
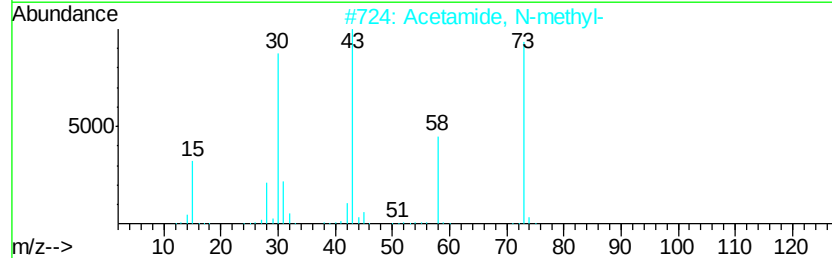
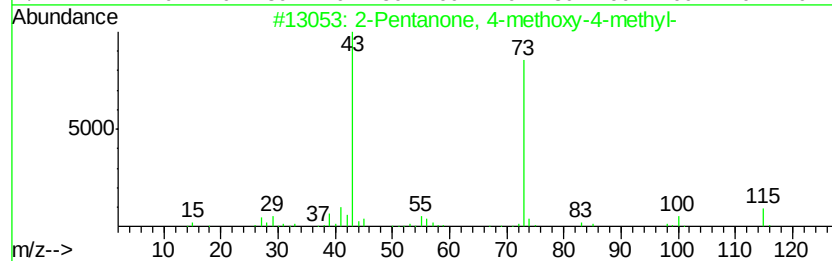
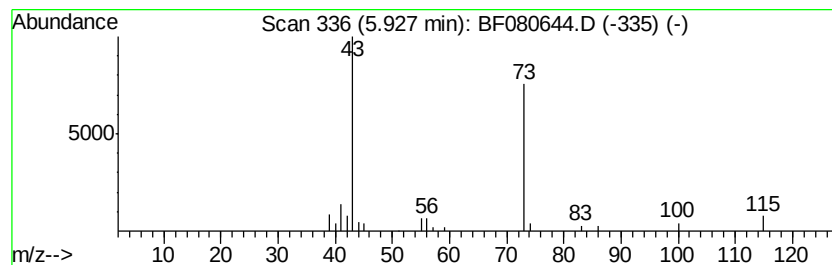
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	2.68 ng	43539	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	25
4		Diacetyl(methoxyimino)methane	143	C6H9NO3	069740-33-6	17
5		N-Formylmorpholine	115	C5H9NO2	004394-85-8	9



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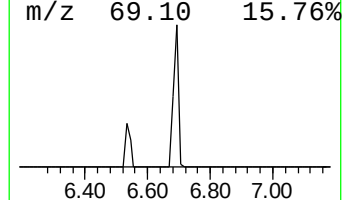
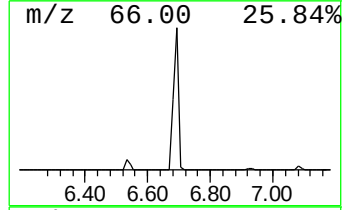
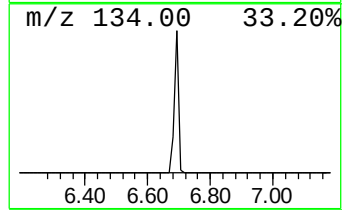
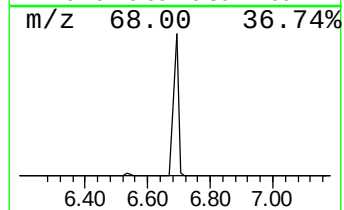
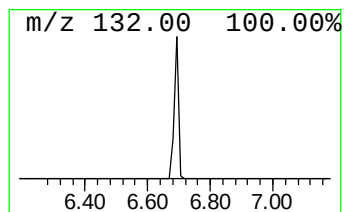
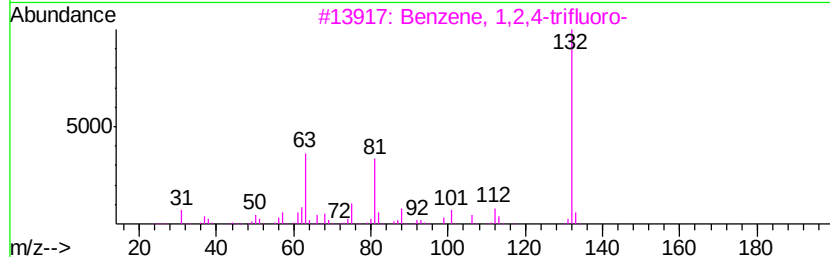
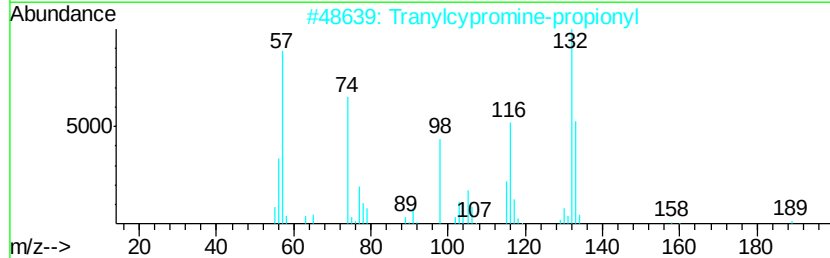
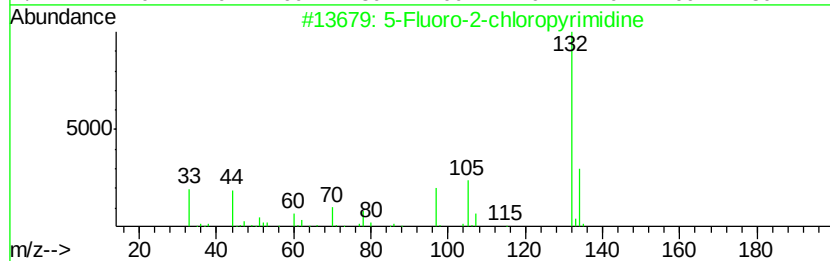
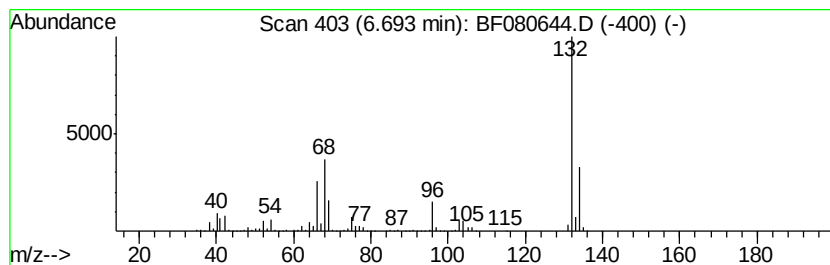
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Peak Number 3 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	72.89 ng	1182240	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.08	13.6	ng	220709	1	6.93	324386	20.0
2-Pentanone, 4-me...	5.93	2.7	ng	43539	1	6.93	324386	20.0
unknown6.69	6.69	72.9	ng	1182240	1	6.93	324386	20.0