

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052215\
 Data File : BF079441.D
 Acq On : 22 May 2015 19:03
 Operator : TP/IZ
 Sample : G2353-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VE2-1

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-BF043015.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	1.324	9	12	15	rVB	260072	367361	8.22%	1.655%
2	1.484	23	26	29	rVB	793642	1105975	24.74%	4.984%
3	1.610	36	37	45	rVV	143792	227824	5.10%	1.027%
4	1.713	45	46	80	rVB	2271100	4470499	100.00%	20.145%
5	4.799	314	316	323	rVB	35727	57003	1.28%	0.257%
6	5.324	360	362	378	rVB	543175	808880	18.09%	3.645%
7	6.627	474	476	485	rBV	463627	506412	11.33%	2.282%
8	6.765	485	488	490	rBV	1513375	1671913	37.40%	7.534%
9	7.050	511	513	516	rBV	416226	387213	8.66%	1.745%
10	7.233	527	529	532	rBV	1618913	1565306	35.01%	7.054%
11	7.748	572	574	577	rBV	1368986	1247437	27.90%	5.621%
12	8.628	649	651	654	rBV	591641	535460	11.98%	2.413%
13	9.976	766	769	771	rBV	2877894	2559654	57.26%	11.534%
14	10.799	838	841	843	rBV	506368	586789	13.13%	2.644%
15	11.771	924	926	929	rBV2	1331435	1664681	37.24%	7.501%
16	12.628	999	1001	1004	rBV	577738	635096	14.21%	2.862%
17	13.371	1064	1066	1070	rBV3	23710	50331	1.13%	0.227%
18	14.628	1173	1176	1178	rBV	2213378	2394594	53.56%	10.791%
19	15.806	1276	1279	1281	rBV2	38517	58151	1.30%	0.262%
20	15.908	1286	1288	1291	rBV	574878	643111	14.39%	2.898%
21	17.646	1437	1440	1444	rBV	502278	647920	14.49%	2.920%

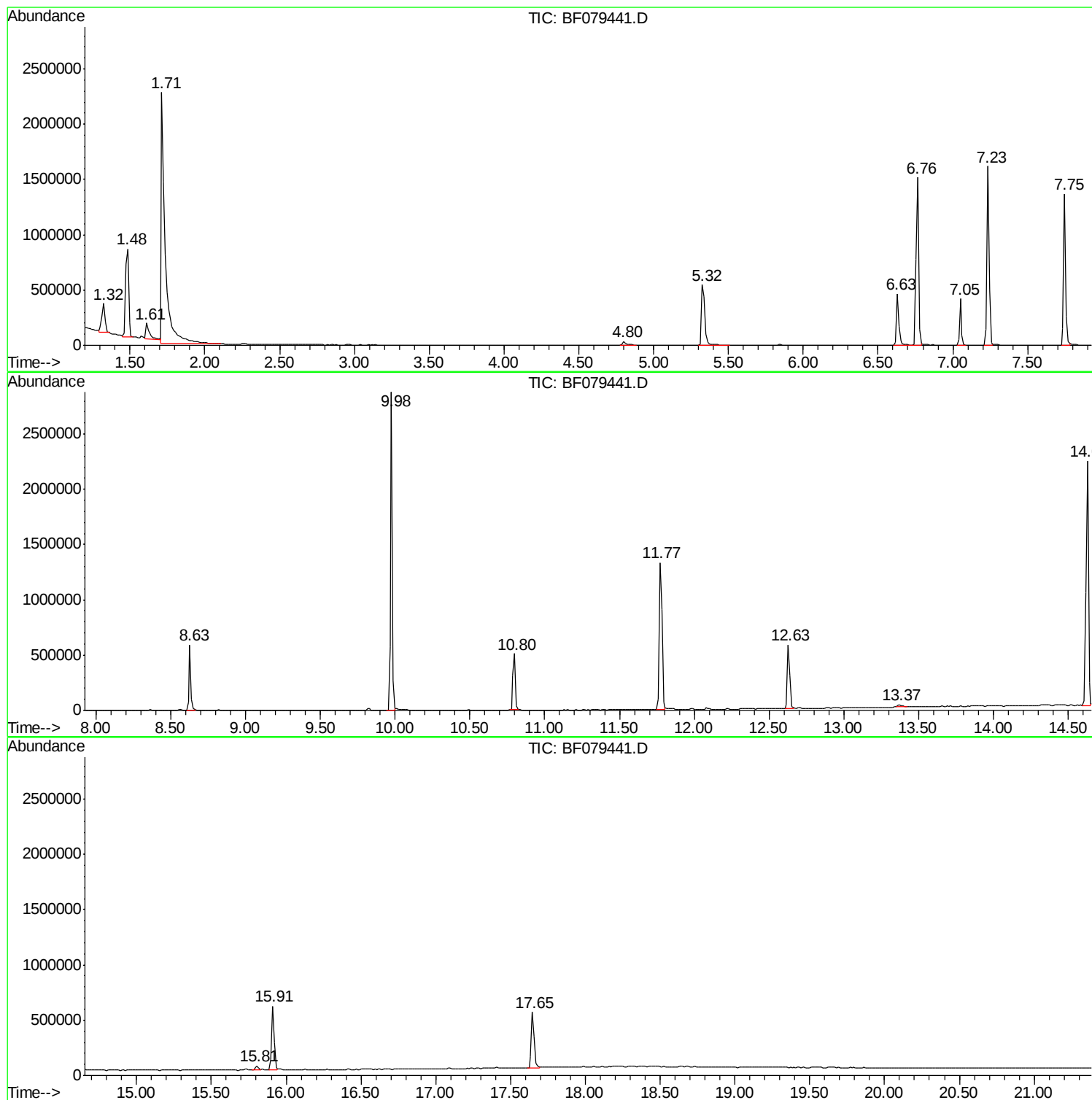
Sum of corrected areas: 22191610

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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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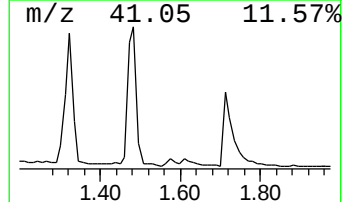
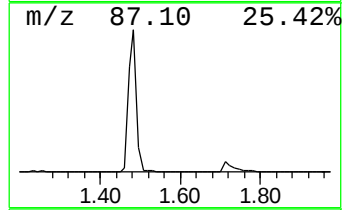
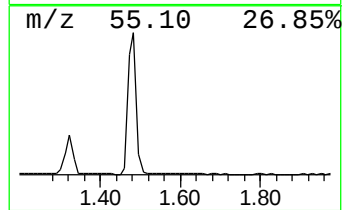
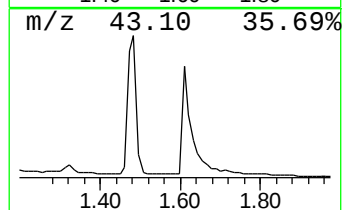
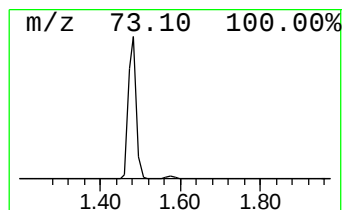
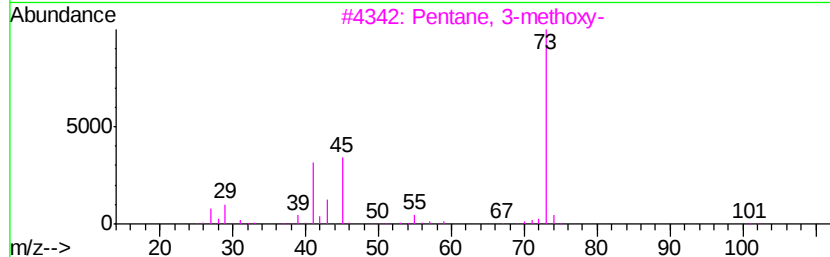
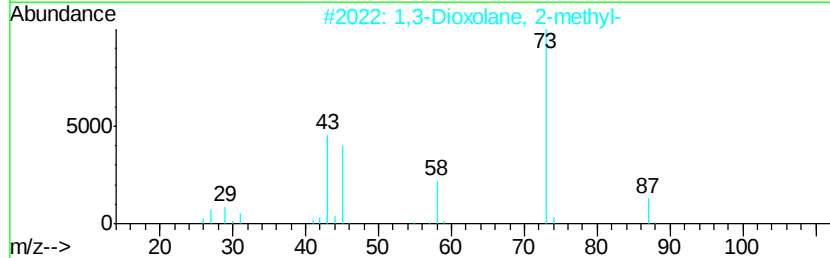
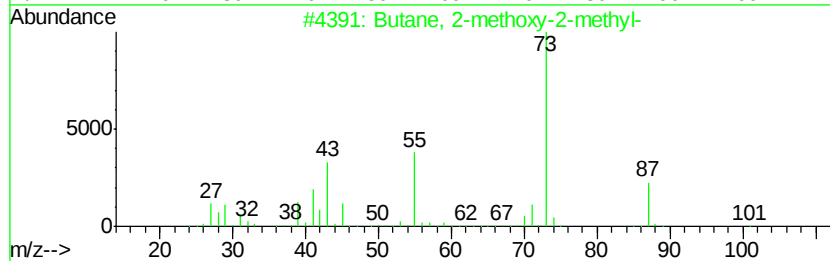
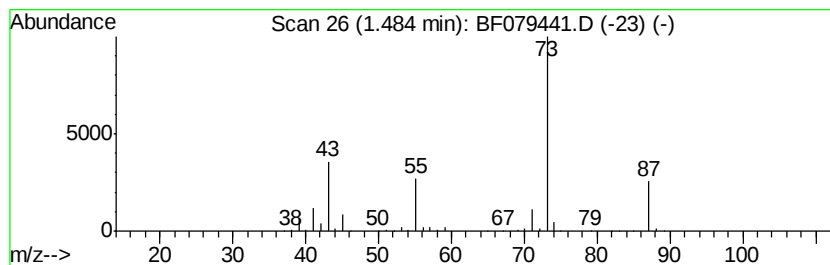
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.48	57.12 ng	1105980	1,4-Dichlorobenzene-d4	7.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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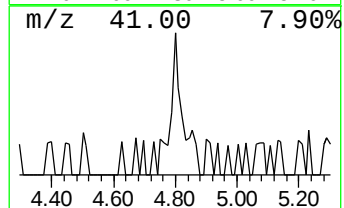
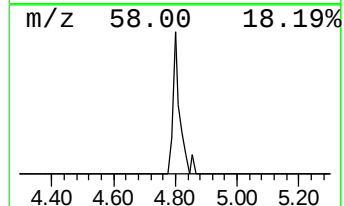
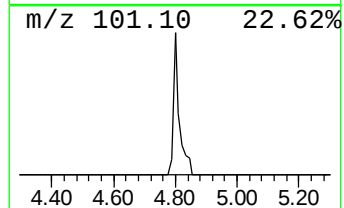
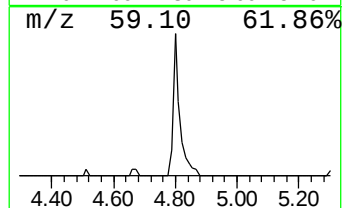
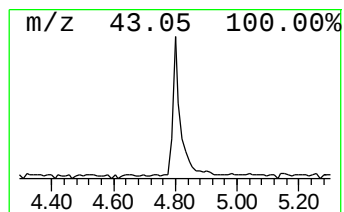
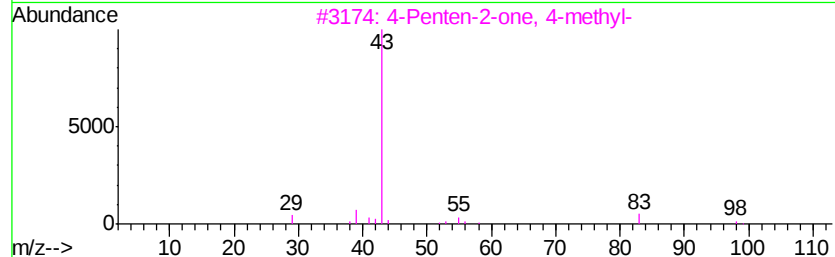
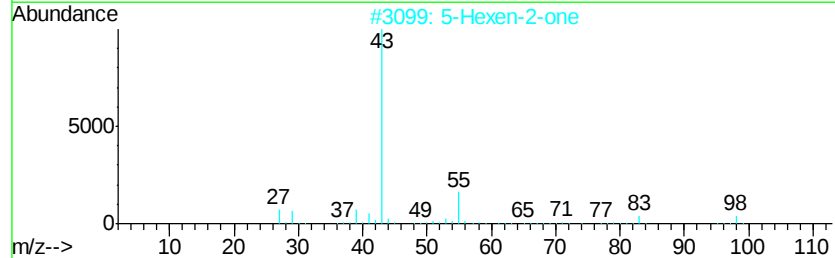
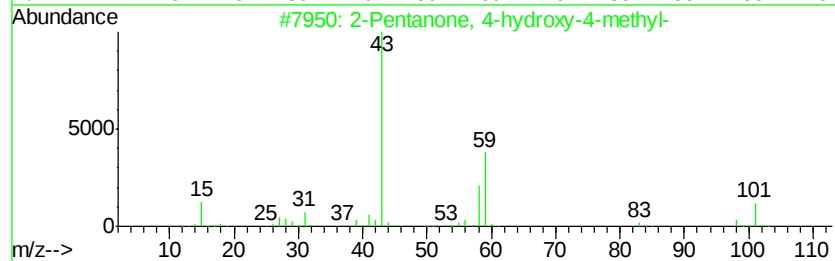
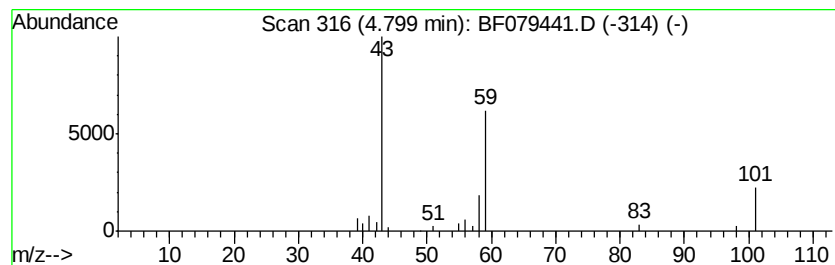
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Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	2.94 ng	57003	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
4		Guanidine	59	CH5N3	000113-00-8	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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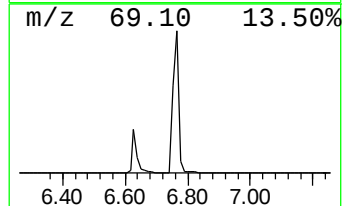
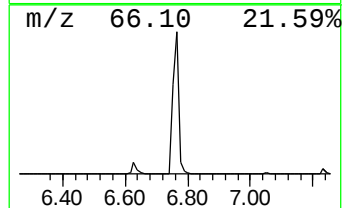
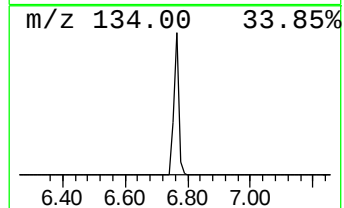
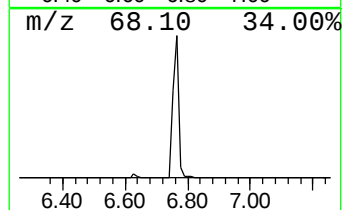
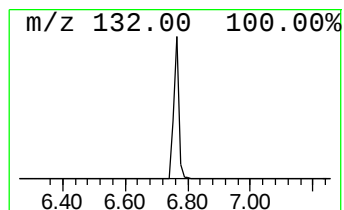
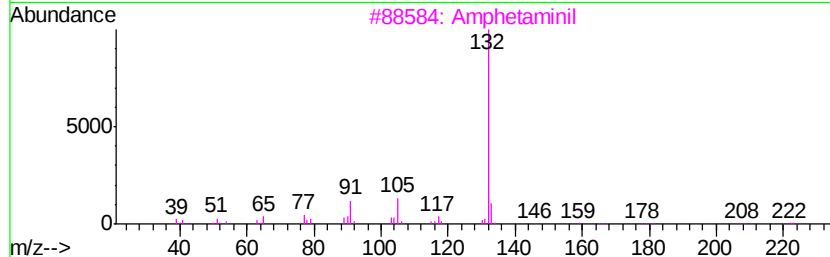
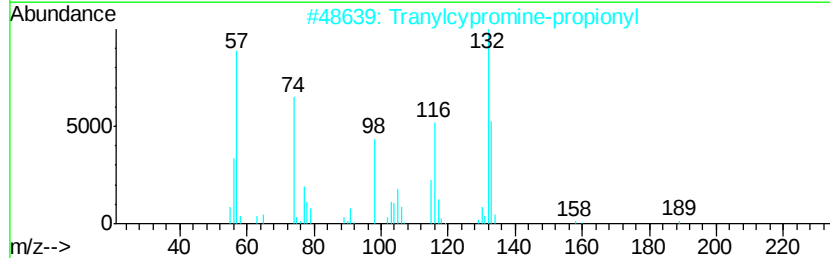
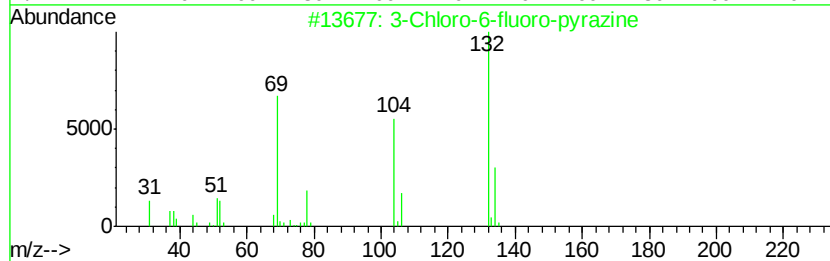
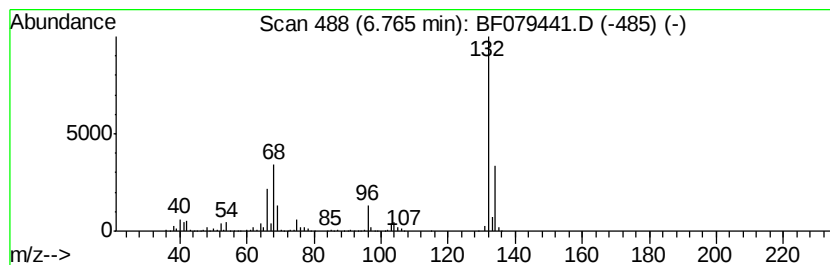
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Peak Number 6 unknown6.76 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	86.36 ng	1671910	1,4-Dichlorobenzene-d4	7.05

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



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.48	57.1	ng	1105980	1	7.05	387213	20.0
2-Pentanone, 4-hy...	4.80	2.9	ng	57003	1	7.05	387213	20.0
unknown6.76	6.76	86.4	ng	1671910	1	7.05	387213	20.0