

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF111117\
 Data File : BF100459.D
 Acq On : 12 Nov 2017 00:51
 Operator : SJ/JU
 Sample : I6369-04
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110617-TIDO-F-U-S

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-BF110817.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	2.210	17	21	28	rVB	79396	110689	2.63%	0.428%
2	5.122	513	516	525	rBV	280855	292521	6.95%	1.132%
3	5.516	579	583	587	rBV	1668627	1484567	35.25%	5.744%
4	6.516	749	753	757	rBV	1183624	1072955	25.48%	4.151%
5	6.669	774	779	782	rBV	2633445	2454873	58.29%	9.498%
6	6.904	815	819	822	rBV	1088939	918348	21.80%	3.553%
7	7.063	841	846	849	rBV	3292060	3145615	74.69%	12.171%
8	7.469	909	915	918	rBV	2524798	2628181	62.40%	10.169%
9	8.186	1032	1037	1040	rBV	1242685	1138595	27.03%	4.405%
10	8.733	1127	1130	1134	rBV	102386	86293	2.05%	0.334%
11	9.263	1214	1220	1223	rBV	4245127	4211711	100.00%	16.295%
12	9.645	1282	1285	1289	rBV	106503	90075	2.14%	0.349%
13	9.939	1329	1335	1343	rVB	1214252	1100650	26.13%	4.259%
14	10.604	1445	1448	1451	rVB	110940	78937	1.87%	0.305%
15	10.651	1452	1456	1459	rBV	273205	230676	5.48%	0.893%
16	10.727	1464	1469	1472	rBV	1514008	1371808	32.57%	5.308%
17	11.427	1584	1588	1591	rBV	970658	883623	20.98%	3.419%
18	13.010	1853	1857	1861	rBV	3040280	3108246	73.80%	12.026%
19	14.063	2032	2036	2040	rBV	950923	830995	19.73%	3.215%
20	15.551	2283	2289	2298	rBV	433805	606564	14.40%	2.347%

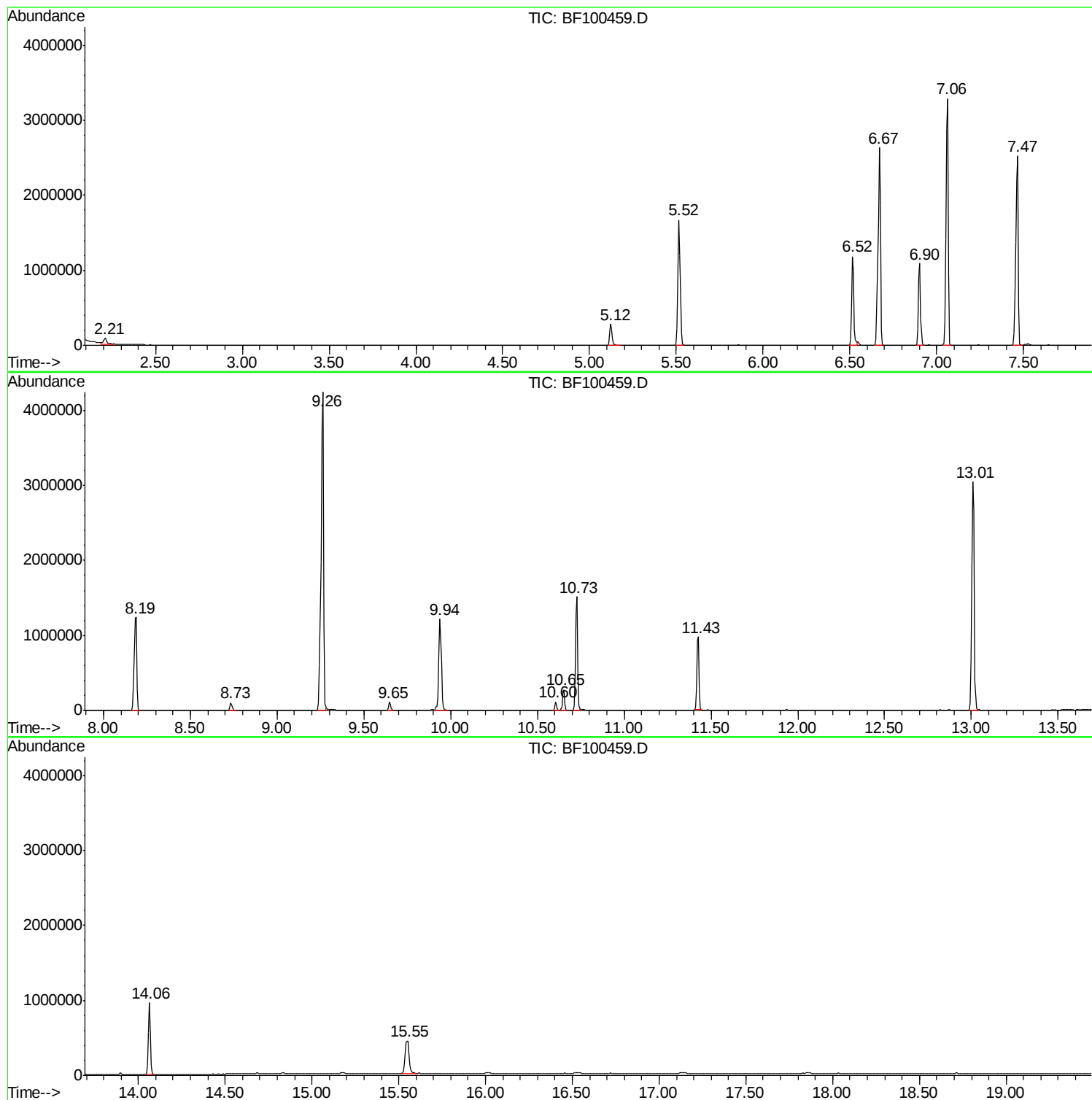
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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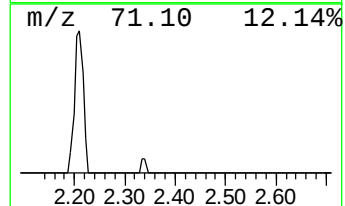
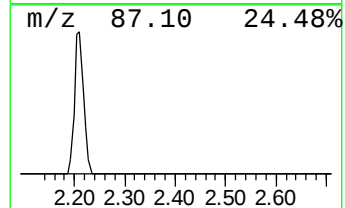
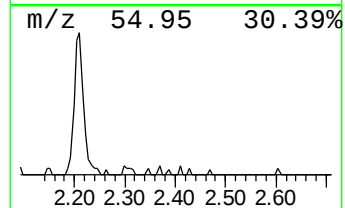
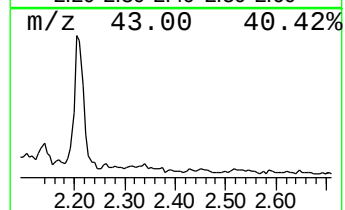
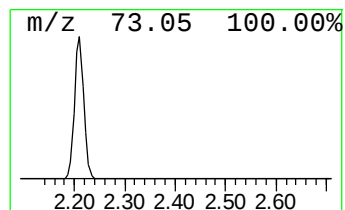
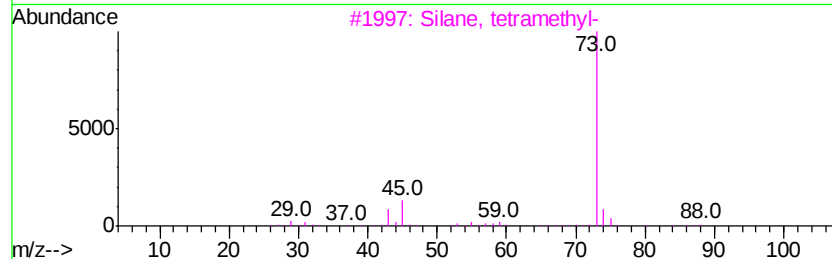
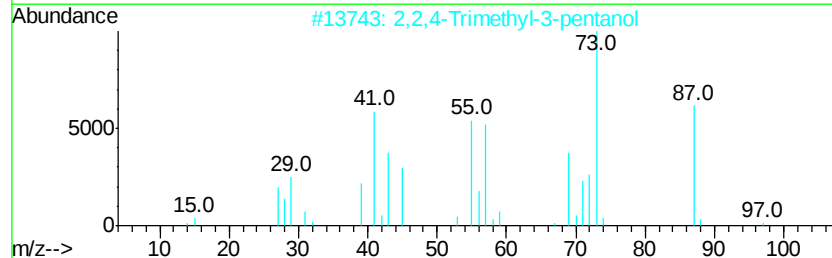
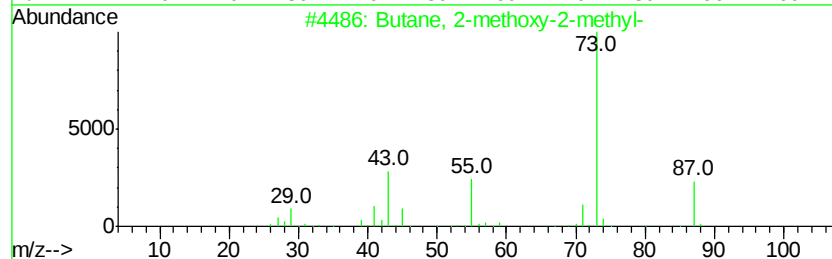
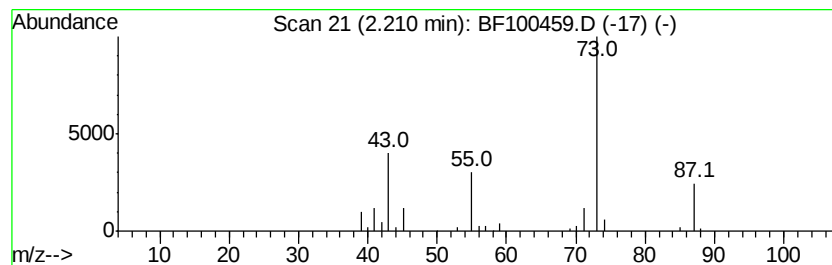
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.21	2.41 ng	110689	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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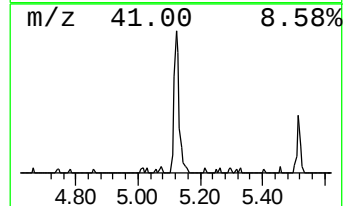
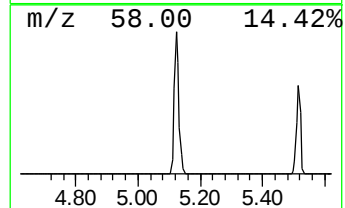
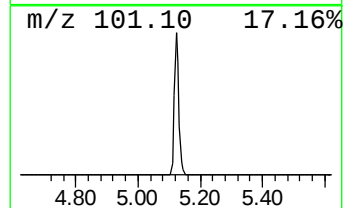
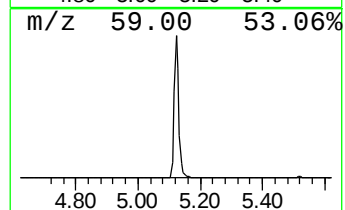
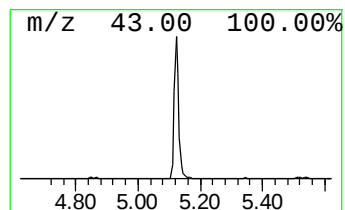
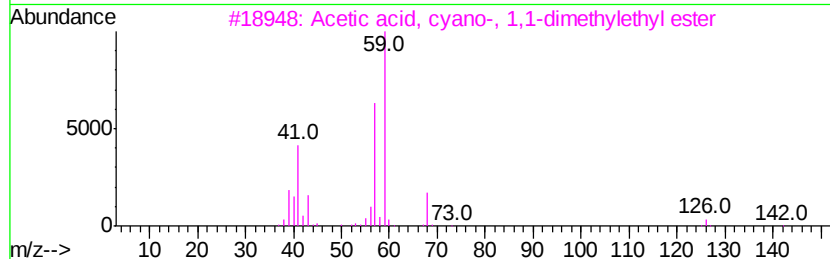
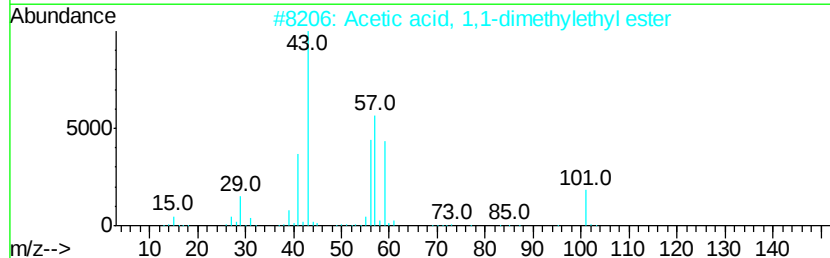
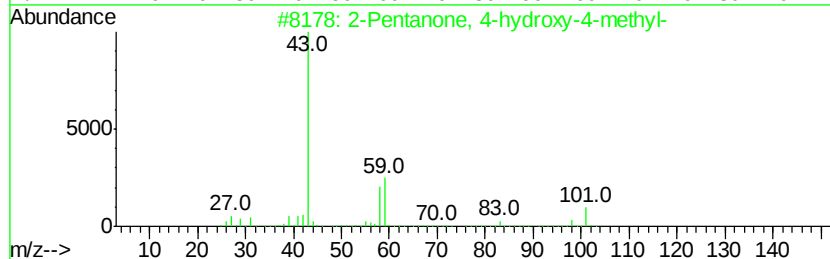
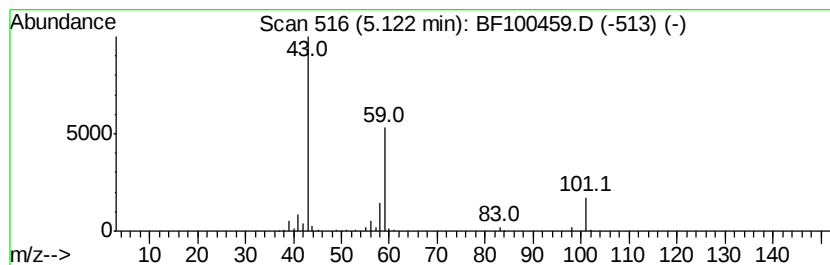
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	6.37 ng	292521	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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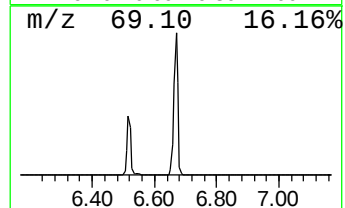
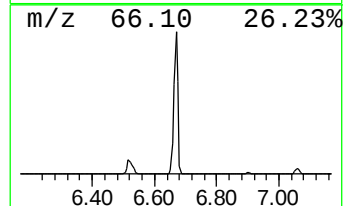
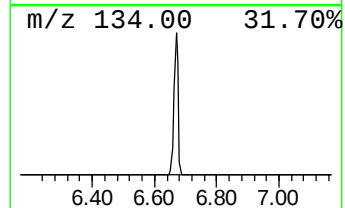
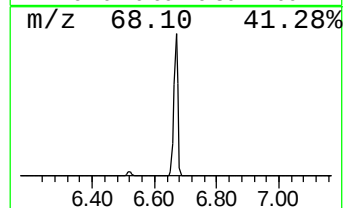
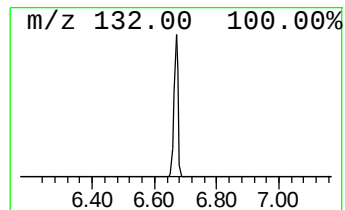
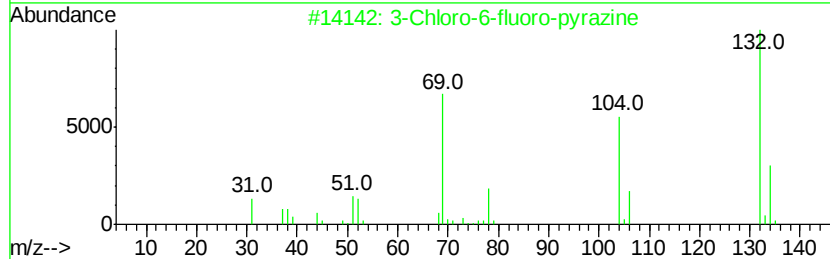
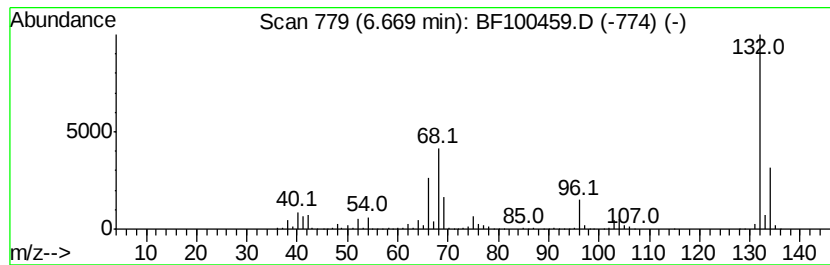
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Peak Number 3 unknown6.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	53.46 ng	2454870	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	17
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
5	1,3,5-Trifluorobenzene			132	C6H3F3	000372-38-3	9



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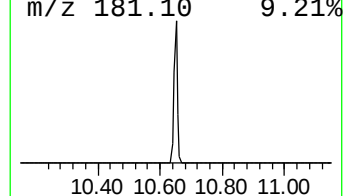
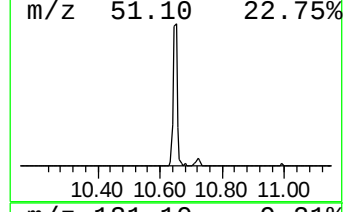
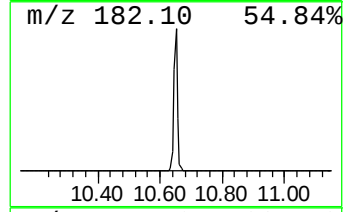
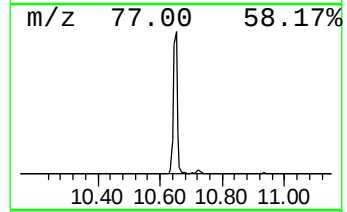
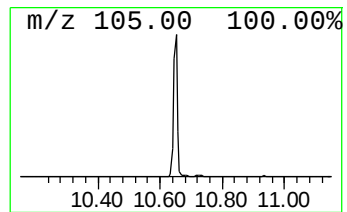
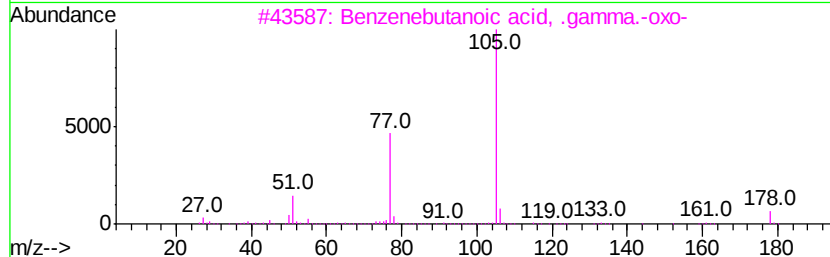
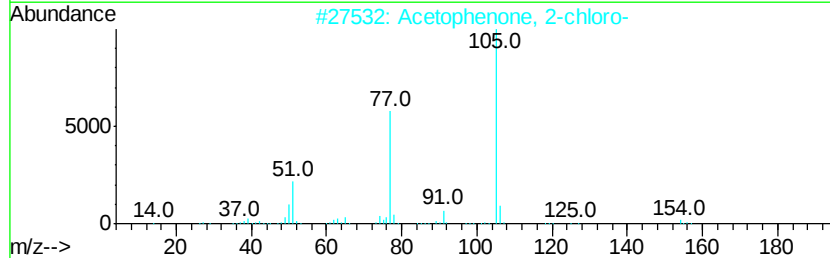
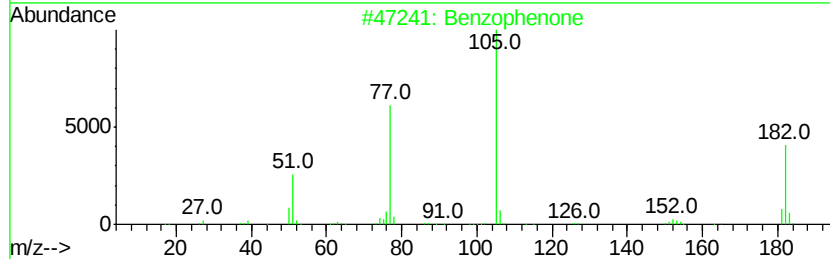
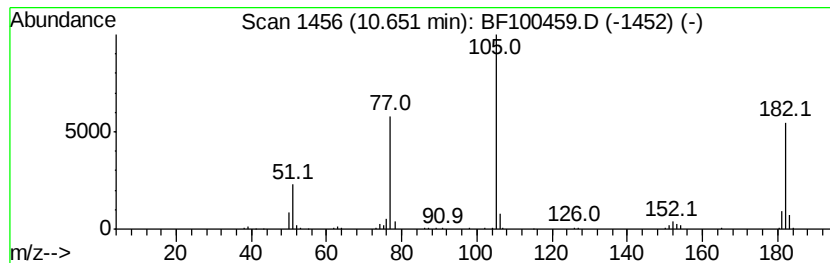
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Peak Number 5 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	4.19 ng	230676	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	97
2		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3		Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
4		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	47
5		Benzoyl chloride	140	C7H5ClO	000098-88-4	47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.21	2.4	ng	110689	1	6.90	918348	20.0
2-Pentanone, 4-hy...	5.12	6.4	ng	292521	1	6.90	918348	20.0
unknown6.67	6.67	53.5	ng	2454870	1	6.90	918348	20.0
Benzophenone	10.65	4.2	ng	230676	3	9.94	1100650	20.0