

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110217\
 Data File : BF100099.D
 Acq On : 2 Nov 2017 21:44
 Operator : SJ/JU
 Sample : I6254-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM-1-2

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-BF102317.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	4.987	490	493	512	rBV	158924	286985	1.01%	0.321%
2	5.393	558	562	598	rBV	1556904	1818046	6.38%	2.033%
3	6.422	733	737	754	rBV	1533119	1836161	6.44%	2.054%
4	6.545	754	758	773	rVB	1989520	1909431	6.70%	2.136%
5	6.769	793	796	806	rBV	540142	489649	1.72%	0.548%
6	6.922	818	822	834	rBV	1633900	1537114	5.39%	1.719%
7	7.339	890	893	910	rBV	1059649	1136226	3.99%	1.271%
8	8.051	1011	1014	1033	rBV	788472	670517	2.35%	0.750%
9	9.128	1192	1197	1200	rBV	2518106	2242542	7.87%	2.508%
10	9.504	1256	1261	1271	rVB	2356277	3295975	11.56%	3.686%
11	9.586	1271	1275	1281	rBV	419364	504413	1.77%	0.564%
12	9.669	1286	1289	1296	rVB2	273651	423057	1.48%	0.473%
13	9.739	1296	1301	1304	rBV2	234809	464541	1.63%	0.520%
14	9.828	1304	1316	1325	rBV5	4484479	28504673	100.00%	31.882%
15	9.986	1340	1343	1345	rBV	1085055	1061141	3.72%	1.187%
16	10.075	1345	1358	1362	rVV4	4137769	20415778	71.62%	22.835%
17	10.145	1366	1370	1377	rVV	4240817	16056171	56.33%	17.958%
18	10.216	1379	1382	1386	rVB2	1196936	1466329	5.14%	1.640%
19	10.251	1386	1388	1396	rVB2	413560	514667	1.81%	0.576%
20	10.433	1415	1419	1426	rVB	367668	373403	1.31%	0.418%
21	10.522	1430	1434	1436	rBV	467420	396467	1.39%	0.443%
22	10.622	1448	1451	1461	rVB	1089009	978536	3.43%	1.094%
23	11.304	1563	1567	1580	rVB	661233	584879	2.05%	0.654%
24	12.880	1830	1835	1838	rBV	1728107	1561785	5.48%	1.747%
25	13.945	2013	2016	2023	rBV	475255	441905	1.55%	0.494%
26	15.404	2260	2264	2275	rVB	338440	436723	1.53%	0.488%

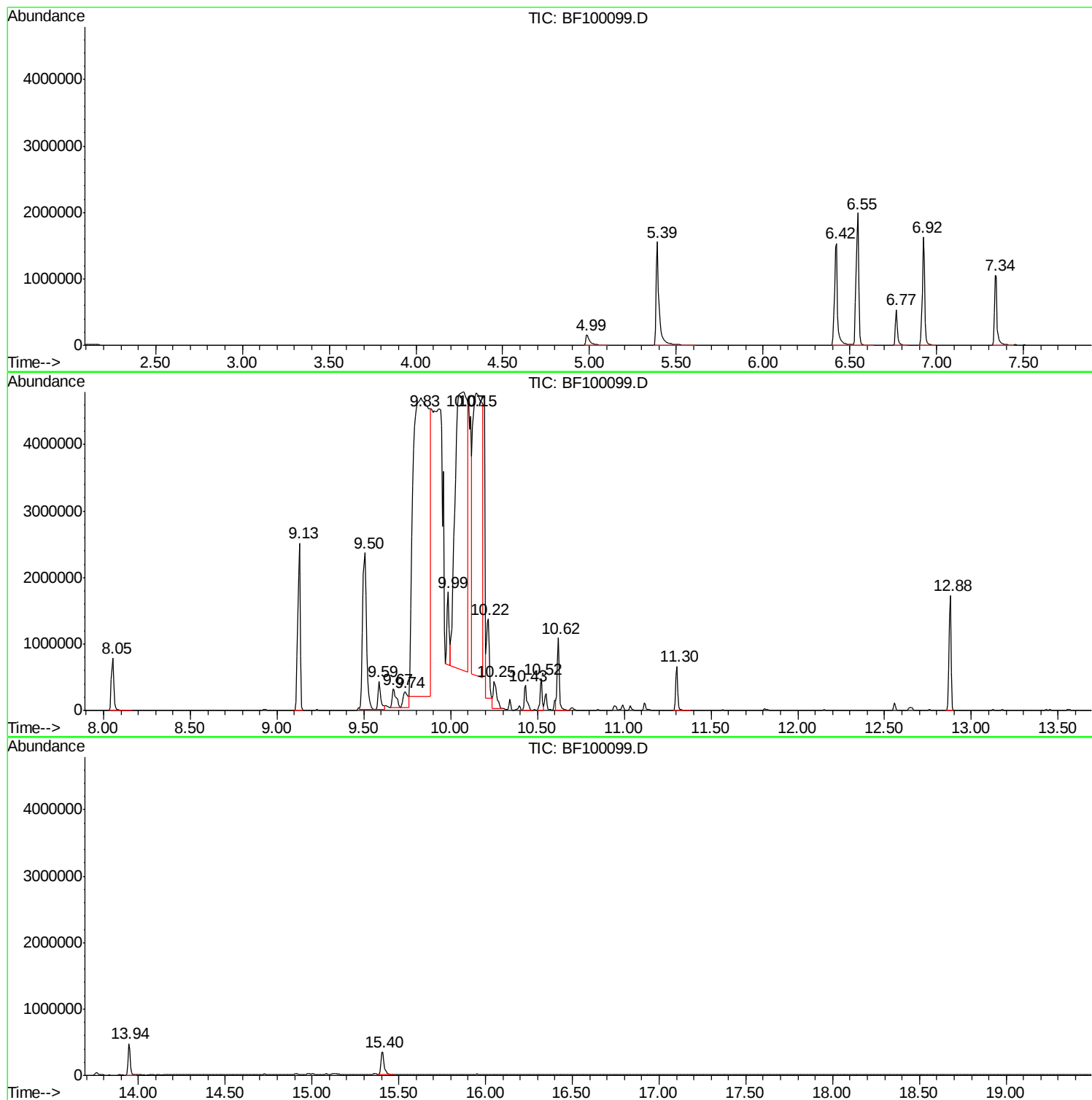
Sum of corrected areas: 89407114

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



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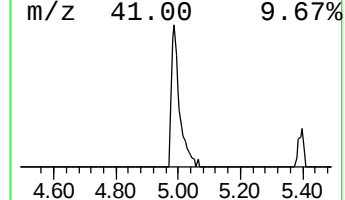
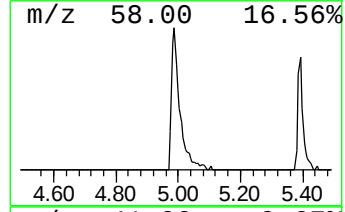
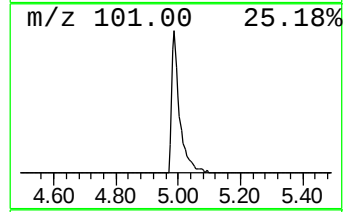
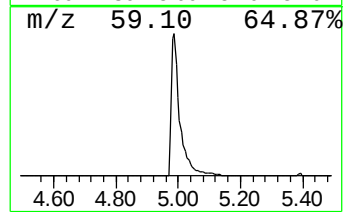
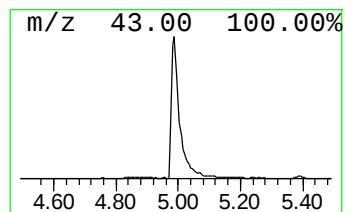
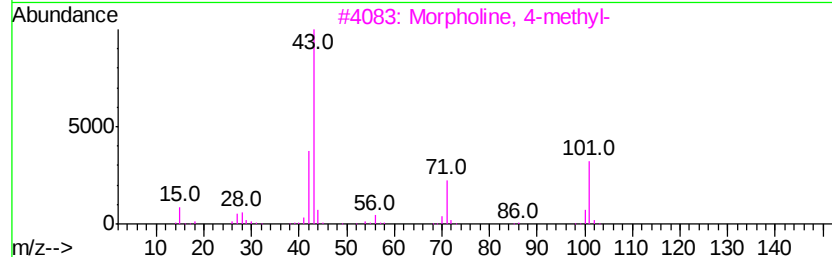
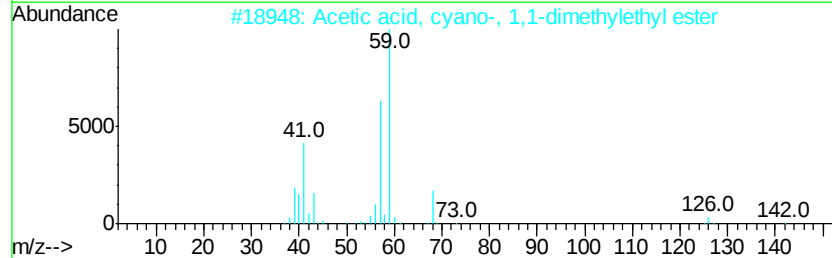
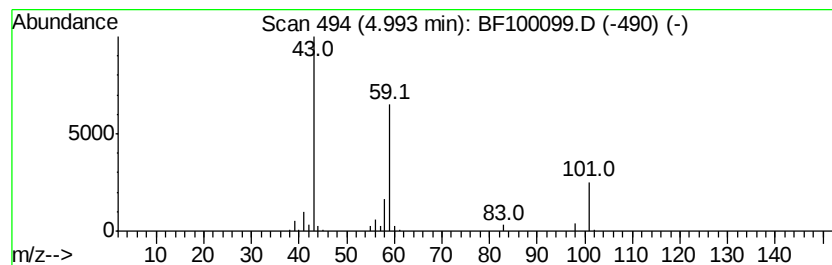
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	11.72 ng	286985	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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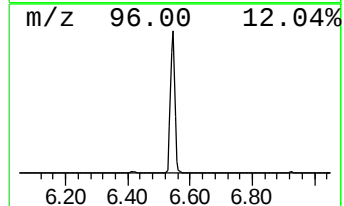
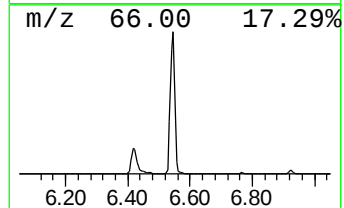
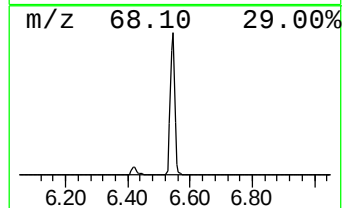
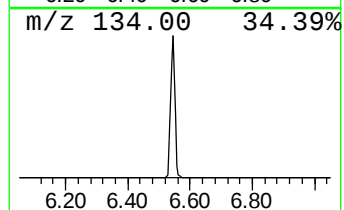
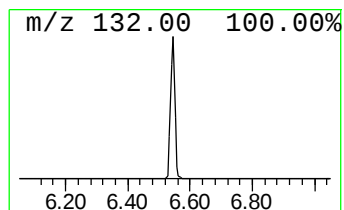
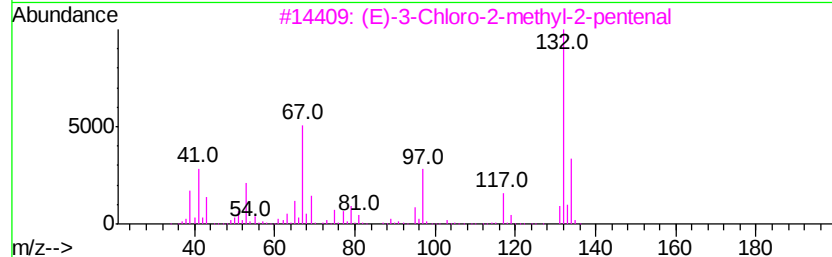
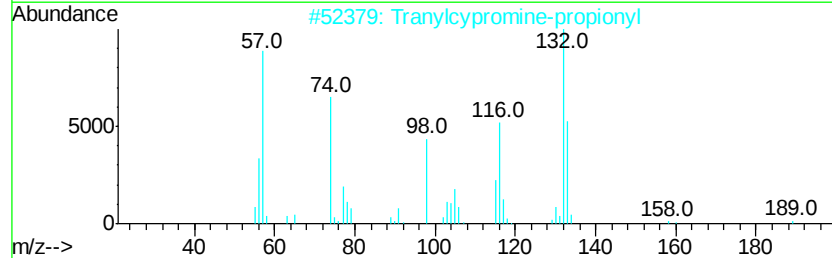
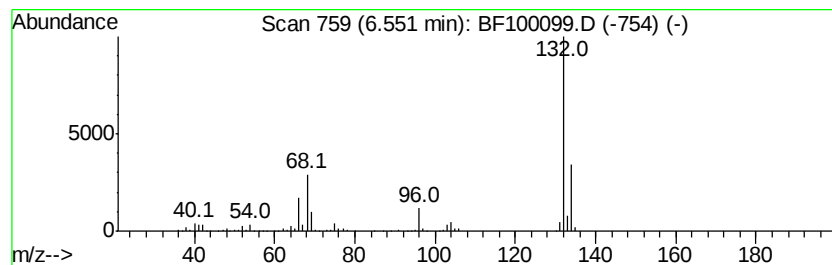
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Peak Number 2 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	77.99 ng	1909430	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		Xenon	132	Xe	007440-63-3	9



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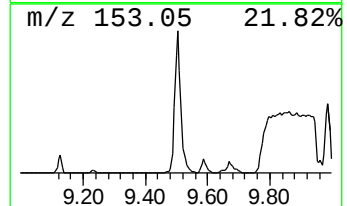
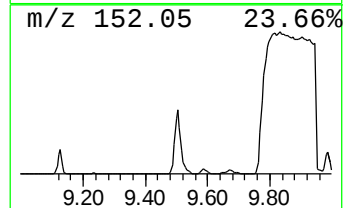
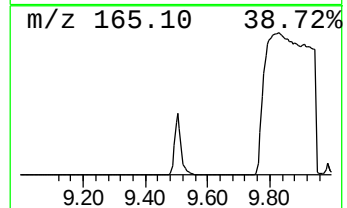
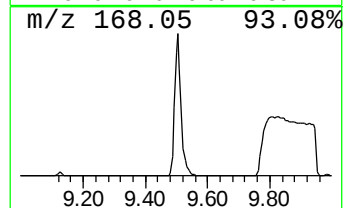
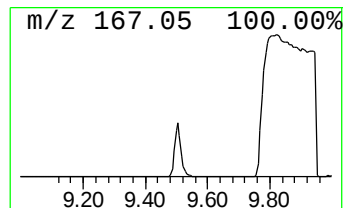
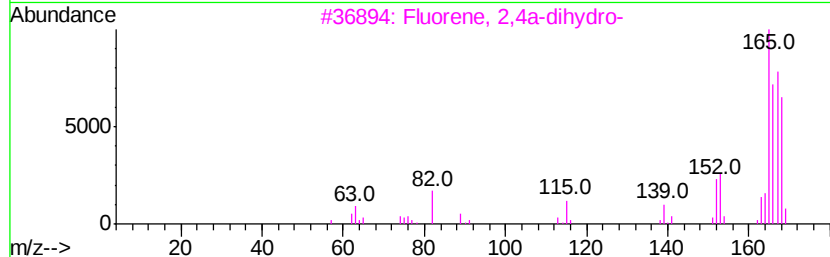
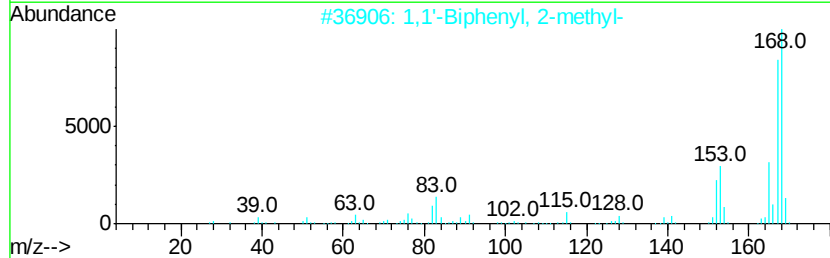
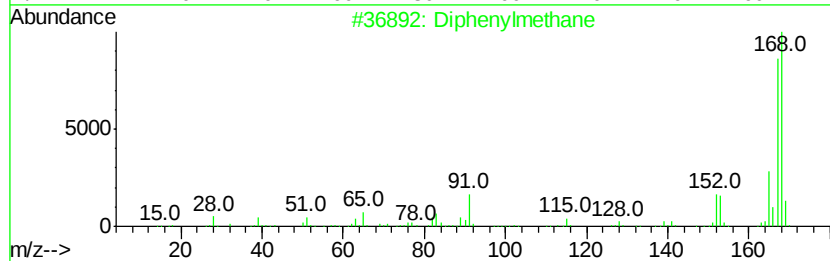
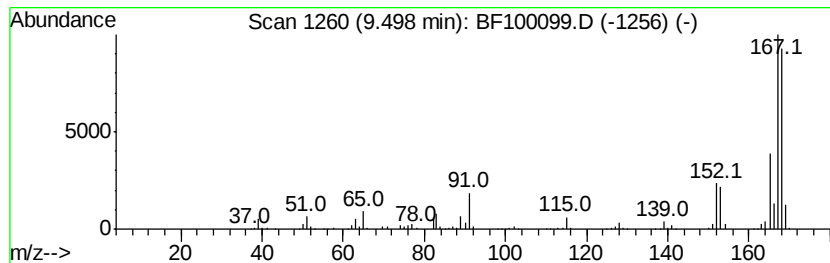
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Peak Number 4 Diphenylmethane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.50	2.31 ng	3295980	Acenaphthene-d10	9.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenylmethane	168	C13H12	000101-81-5	93
2	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	81
3	Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	64
4	Nordiphenamid	225	C15H15NO	000954-21-2	64
5	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	60



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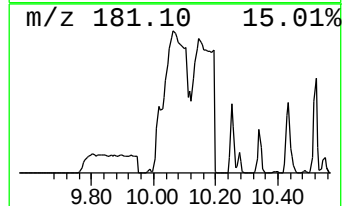
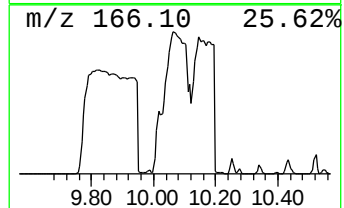
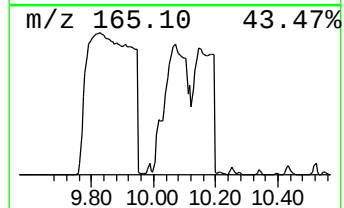
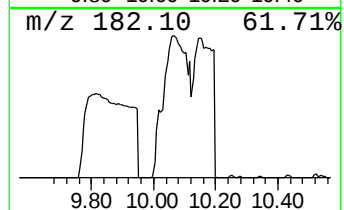
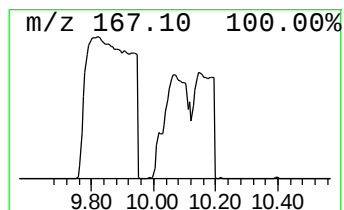
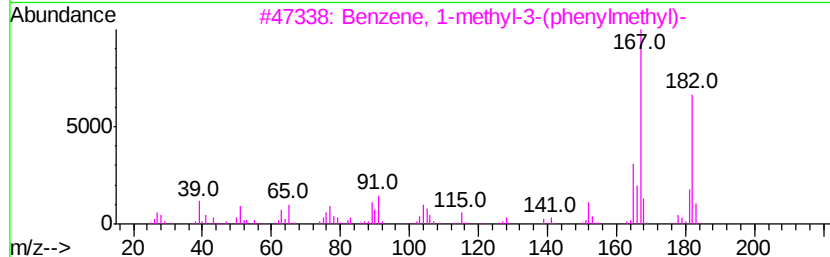
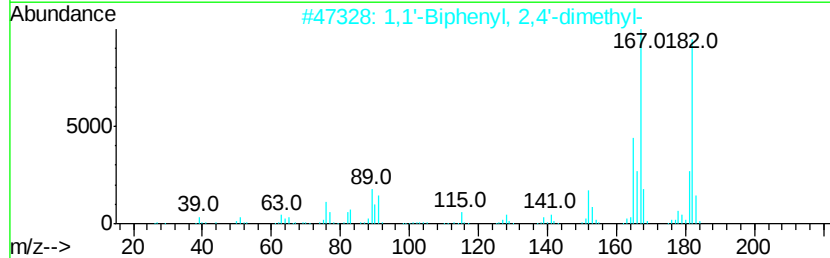
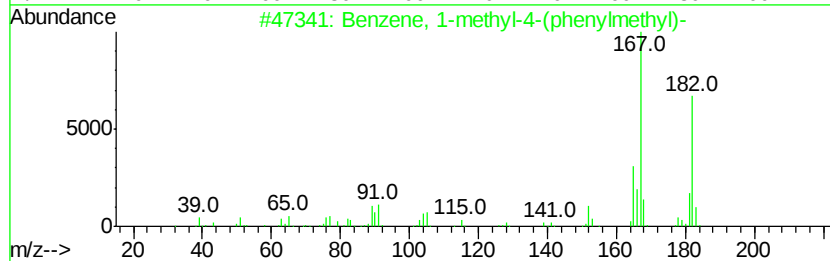
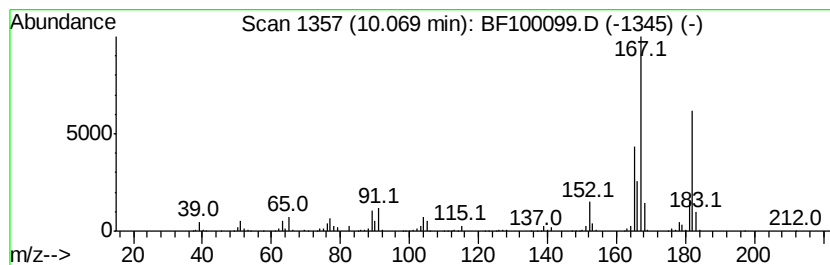
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Peak Number 5 Benzene, 1-methyl-4-(phenyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.07	14.32 ng	20415800	Acenaphthene-d10	9.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-4-(phenylmethyl)-	182 C14H14	000620-83-7 96
2		1,1'-Biphenyl, 2,4'-dimethyl-	182 C14H14	000611-61-0 94
3		Benzene, 1-methyl-3-(phenylmethyl)-	182 C14H14	000620-47-3 93
4		2,2'-Dimethylbiphenyl	182 C14H14	000605-39-0 93
5		Benzene, 1-methyl-2-(phenylmethyl)-	182 C14H14	000713-36-0 80



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	11.7	ng	286985	1	6.77	489649	20.0
unknown6.55	6.55	78.0	ng	1909430	1	6.77	489649	20.0
Diphenylmethane	9.50	2.3	ng	3295980	3	9.83	28504700	20.0
Benzene, 1-methyl...	10.07	14.3	ng	20415800	3	9.83	28504700	20.0