

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101119\
 Data File : BP000639.D
 Acq On : 11 Oct 2019 18:22
 Operator : JU
 Sample : K5266-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AMENDED-COMPOSITE-A

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP100119.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	4.993	481	494	497	rBV	4749006	11273652	100.00%	21.199%
2	5.369	555	558	576	rVB	856700	843995	7.49%	1.587%
3	6.381	727	730	742	rBV2	2030833	2146150	19.04%	4.036%
4	6.516	750	753	761	rVB	1504538	1174961	10.42%	2.209%
5	6.746	789	792	796	rBV	1748032	1442923	12.80%	2.713%
6	6.904	815	819	822	rBV	5400825	4466579	39.62%	8.399%
7	7.310	883	888	891	rBV	3637598	3080029	27.32%	5.792%
8	8.034	1007	1011	1014	rBV	2208765	1864565	16.54%	3.506%
9	9.116	1191	1195	1198	rBV	7501185	6125599	54.34%	11.519%
10	9.510	1258	1262	1265	rBV	1213135	893018	7.92%	1.679%
11	9.792	1307	1310	1314	rVB	2601586	2359912	20.93%	4.438%
12	10.587	1442	1445	1450	rBV	784599	650627	5.77%	1.223%
13	11.151	1538	1541	1555	rBV2	264424	357287	3.17%	0.672%
14	11.292	1561	1565	1568	rBV	3207585	2599533	23.06%	4.888%
15	11.363	1574	1577	1581	rVB2	222369	193047	1.71%	0.363%
16	12.516	1770	1773	1776	rVB	173764	148091	1.31%	0.278%
17	12.745	1809	1812	1816	rVB	176073	176287	1.56%	0.331%
18	12.898	1834	1838	1841	rBV	7311932	6951616	61.66%	13.072%
19	13.822	1992	1995	2003	rBV2	181351	252168	2.24%	0.474%
20	13.963	2015	2019	2022	rBV	2738172	2620362	23.24%	4.927%
21	14.457	2100	2103	2106	rBV	148674	137851	1.22%	0.259%
22	14.763	2152	2155	2165	rVB2	165521	193312	1.71%	0.364%
23	15.104	2209	2213	2219	rVB2	193898	230148	2.04%	0.433%
24	15.375	2256	2259	2264	rVV	85574	117905	1.05%	0.222%
25	15.439	2265	2270	2274	rVV	1991905	2417555	21.44%	4.546%
26	15.486	2275	2278	2283	rVB	120529	159959	1.42%	0.301%
27	15.927	2349	2353	2357	rBV	92834	140884	1.25%	0.265%
28	16.151	2387	2391	2404	rBV10	51213	161660	1.43%	0.304%

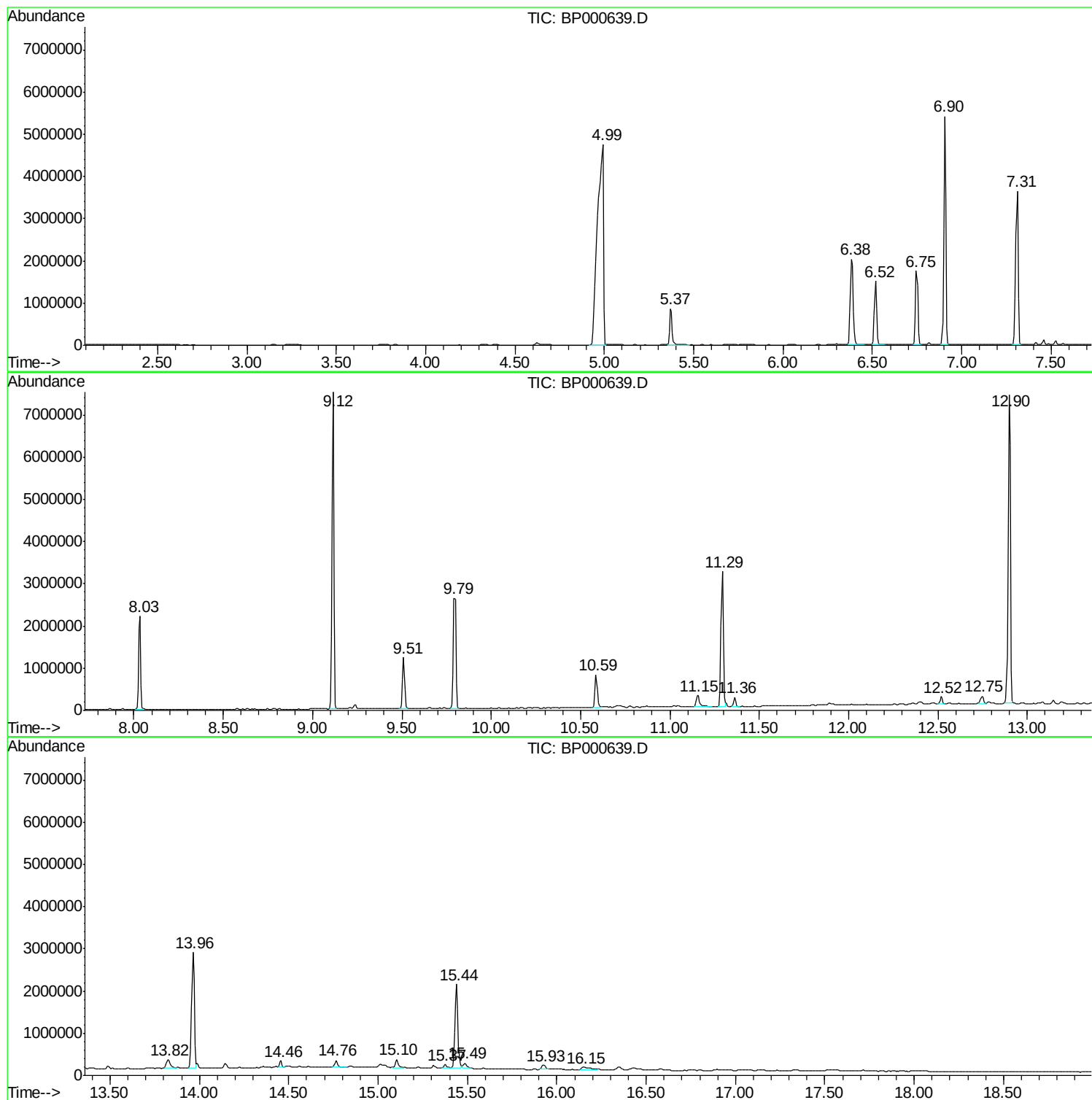
Sum of corrected areas: 53179675

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.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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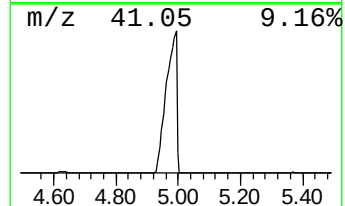
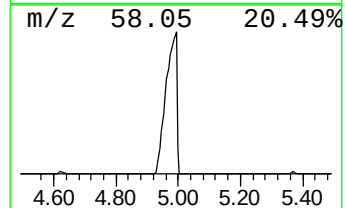
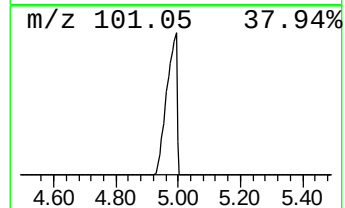
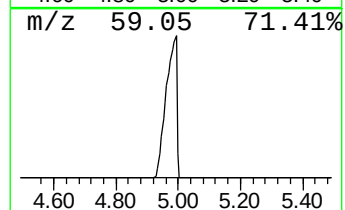
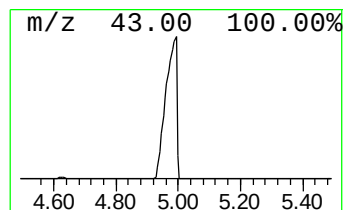
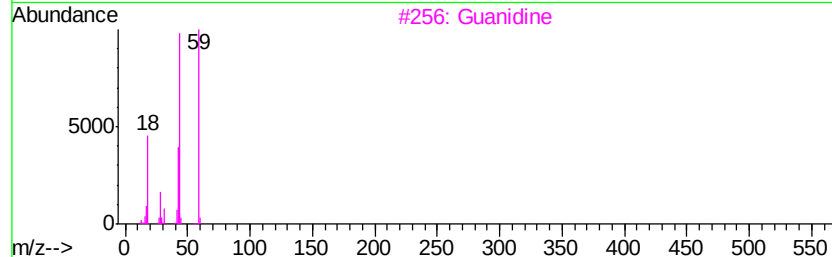
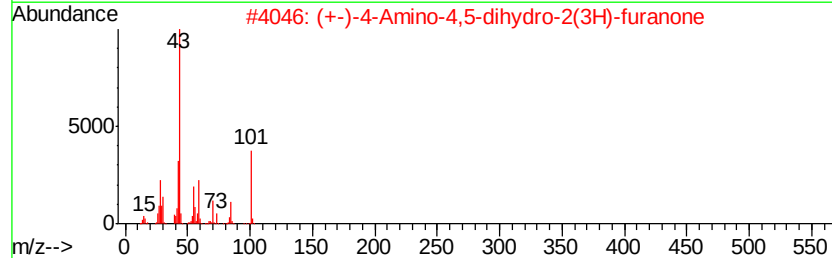
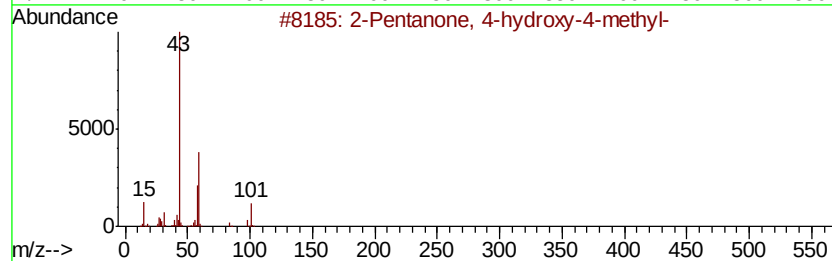
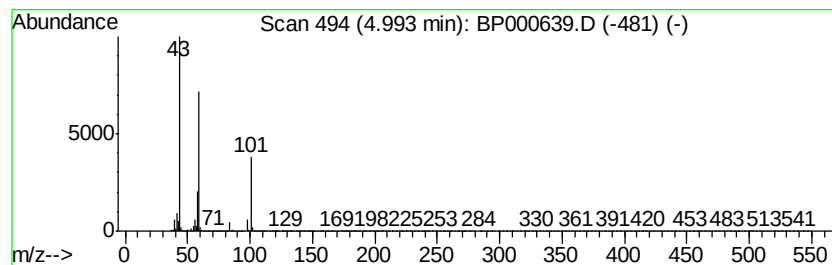
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.99 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	156.26 ng	11273700	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	43
3			Guanidine	59	CH5N3	000113-00-8	38
4			2-Nonanone	142	C9H18O	000821-55-6	12
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	12



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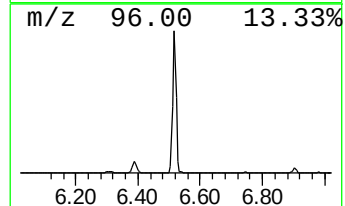
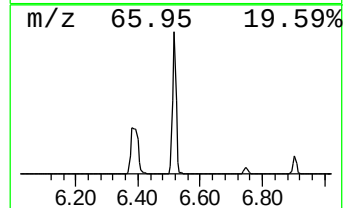
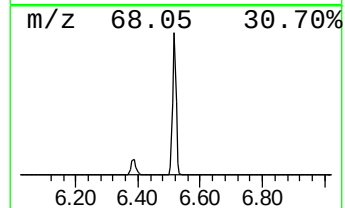
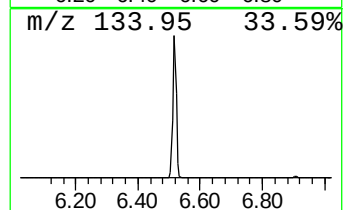
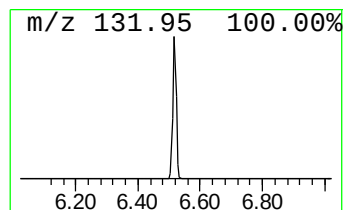
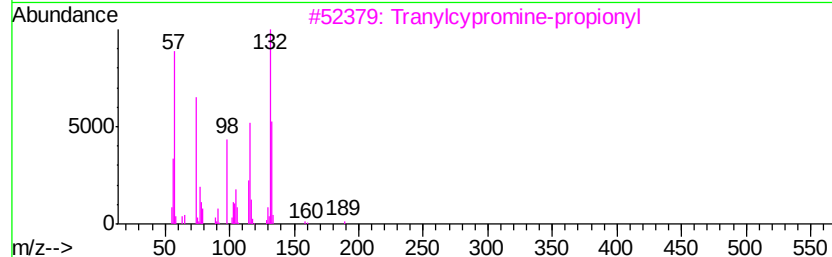
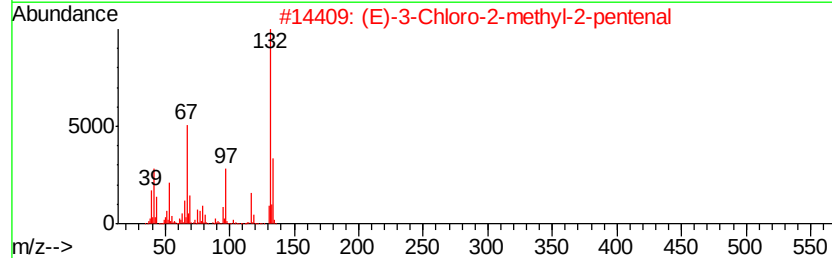
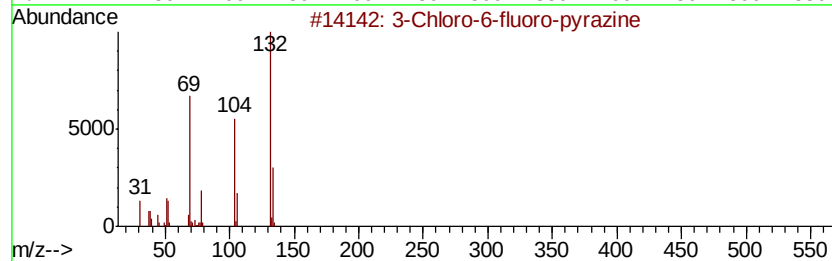
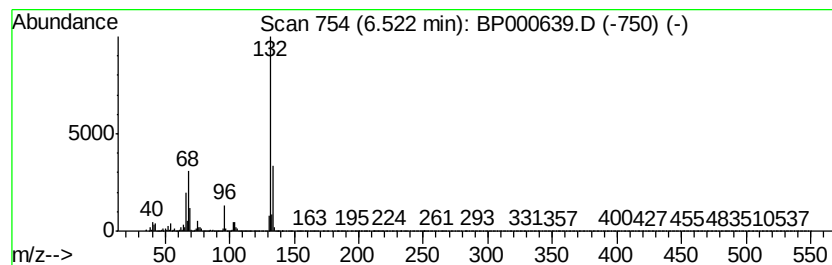
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Peak Number 2 unknown6.52 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	16.29 ng	1174960	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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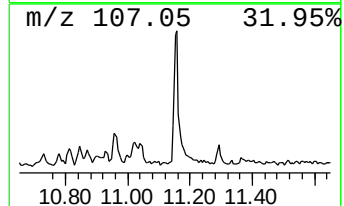
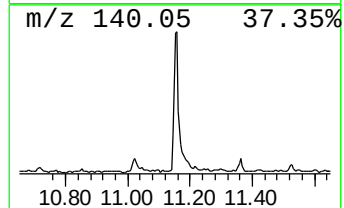
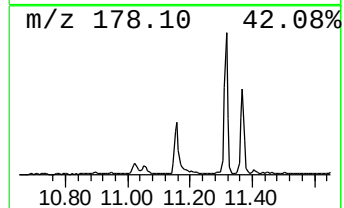
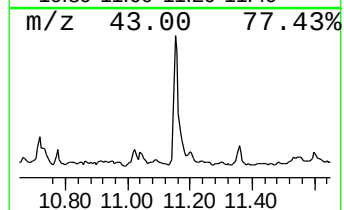
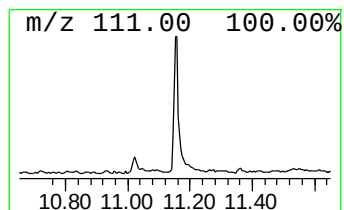
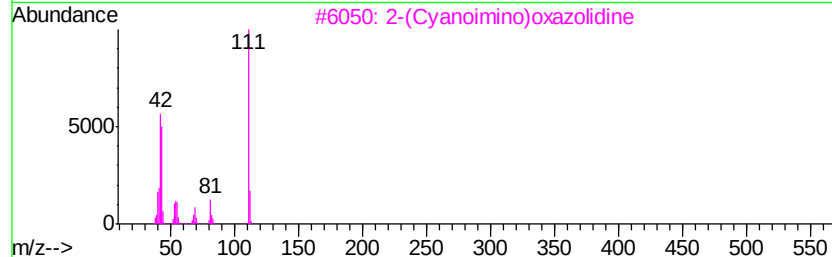
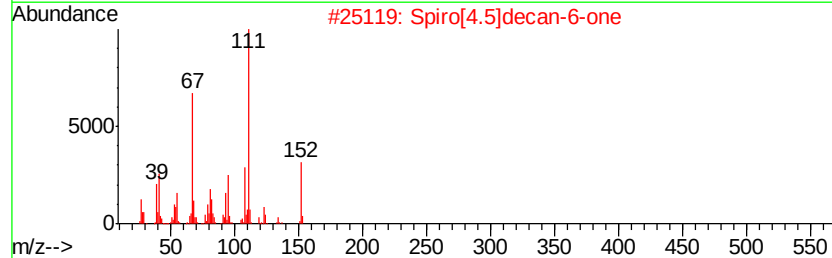
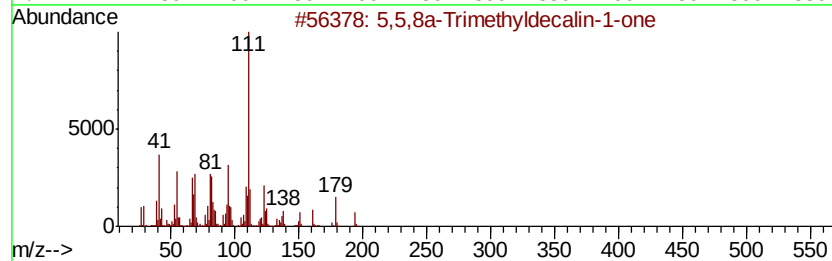
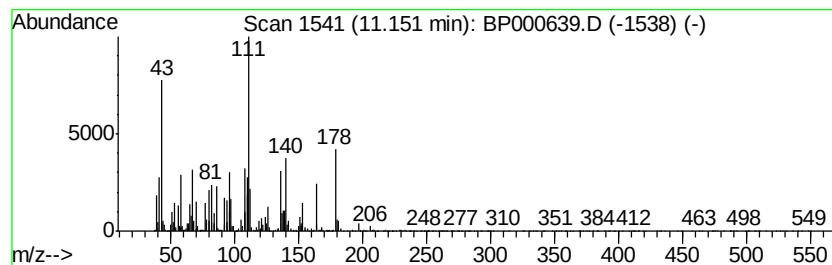
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Peak Number 4 unknown11.15 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	2.75 ng	357287	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,5,8a-Trimethyldecalin-1-one	194	C13H22O	1000195-96-1	32
2		Spiro[4.5]decan-6-one	152	C10H16O	013388-94-8	30
3		2-(Cyanoimino)oxazolidine	111	C4H5N3O	050411-06-8	25
4		2-Methylcyclohexan-1-on-3-ylbuta...	362	C17H30O6S	344552-36-9	25
5		1-Methyl-3-n-propyl-2-pyrazolin-...	140	C7H12N2O	031272-04-5	22



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown4.99	4.99	156.3	ng	11273700	1	6.75	1442920	20.0
unknown6.52	6.52	16.3	ng	1174960	1	6.75	1442920	20.0
unknown11.15	11.15	2.8	ng	357287	4	11.29	2599530	20.0