

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\
 Data File : BM033370.D
 Acq On : 09 Dec 2021 22:48
 Operator : CG/JU
 Sample : M4960-11DL2 300X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGKS4DL2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_M\METHODS\SFAM-EPA-BM120921.M
 Title : SVOA CALIBRATION

Signal : TIC: BM033370.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.555	390	394	399	rVB5	5219	8623	0.23%	0.116%
2	6.596	567	571	576	rVB4	5344	8833	0.23%	0.119%
3	7.907	788	794	807	rBV	195958	328890	8.61%	4.415%
4	8.819	944	949	960	rBV5	6543	15008	0.39%	0.201%
5	9.743	1100	1106	1113	rBV6	5113	9942	0.26%	0.133%
6	10.701	1262	1269	1276	rBV	249681	429897	11.26%	5.770%
7	10.748	1276	1277	1284	rVB4	4725	5708	0.15%	0.077%
8	10.872	1284	1298	1313	rBV	2198183	3819576	100.00%	51.270%
9	12.243	1525	1531	1543	rBV	158242	275527	7.21%	3.698%
10	14.531	1913	1920	1928	rBV	346445	537763	14.08%	7.218%
11	15.336	2049	2057	2072	rBV2	23934	70587	1.85%	0.947%
12	17.272	2379	2386	2399	rBV	420152	611498	16.01%	8.208%
13	21.430	3088	3093	3102	rBV	476349	665641	17.43%	8.935%
14	23.754	3482	3488	3499	rVB2	323616	662454	17.34%	8.892%

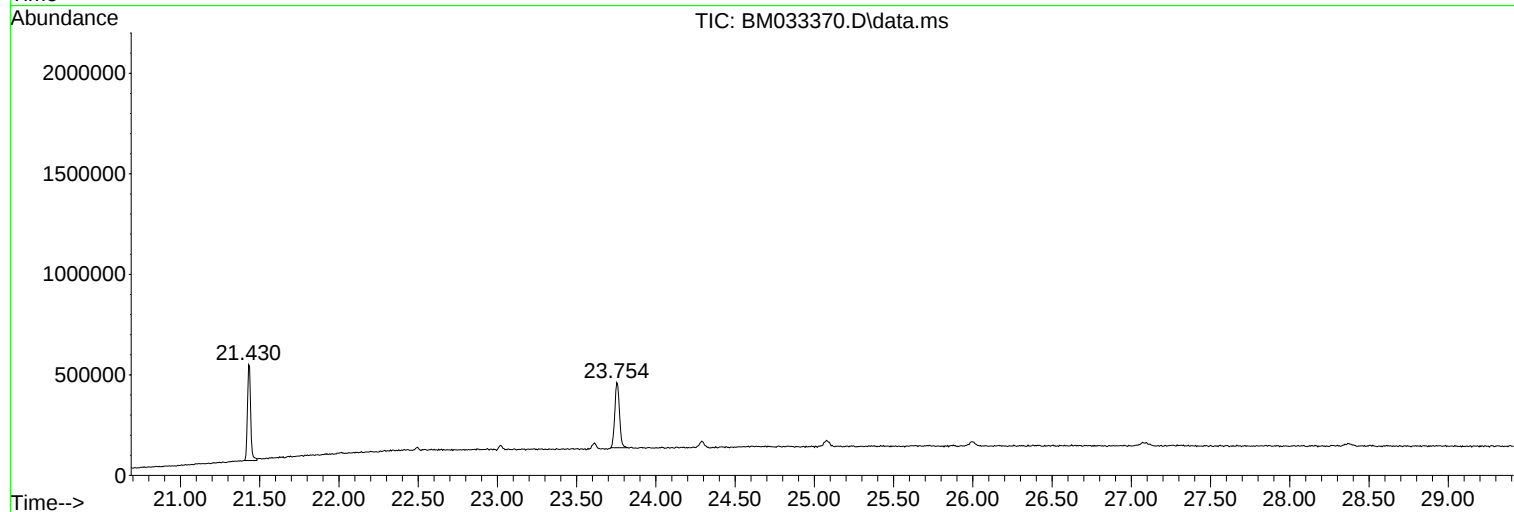
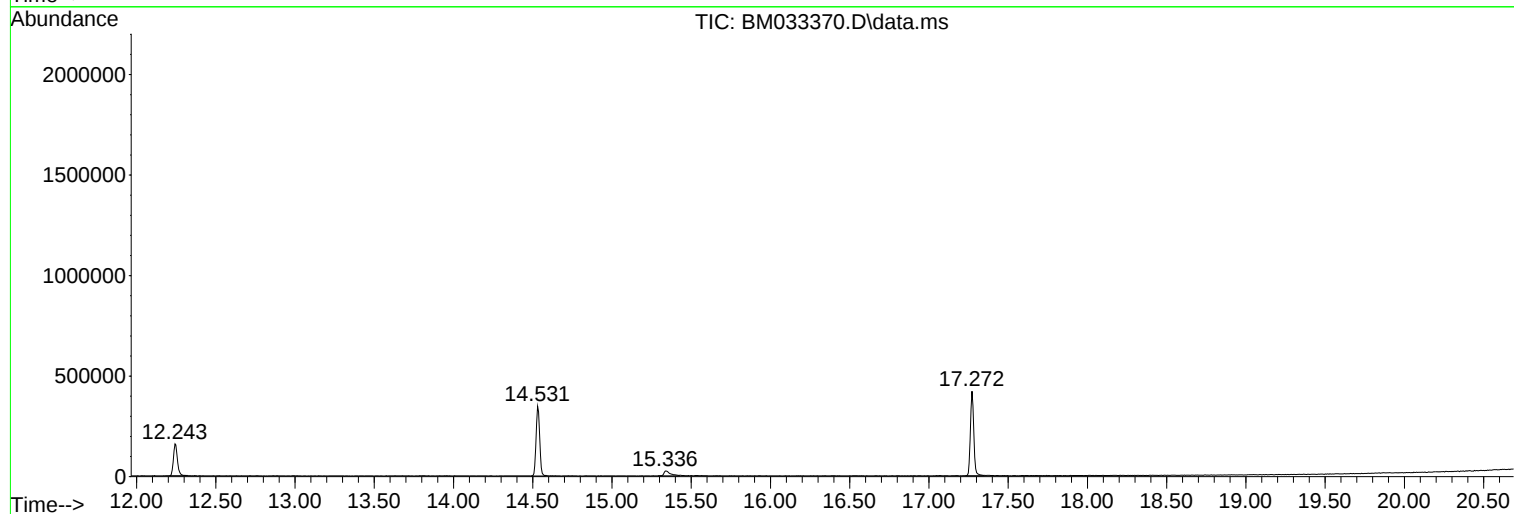
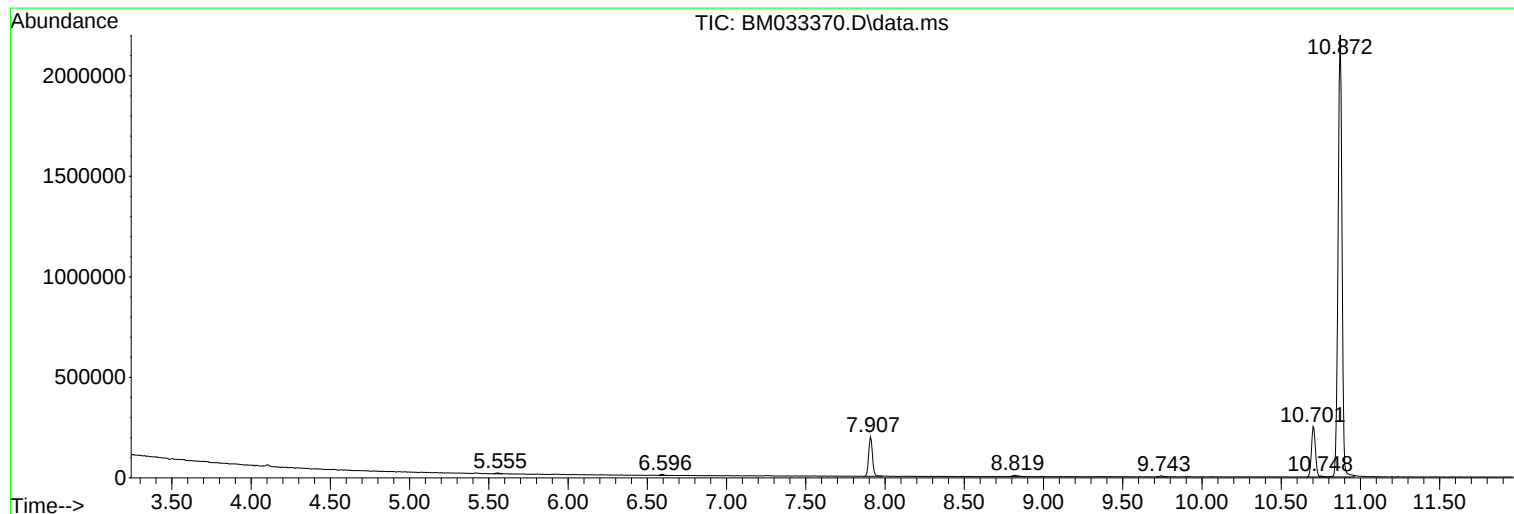
Sum of corrected areas: 7449947

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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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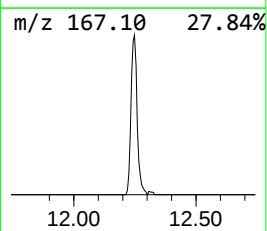
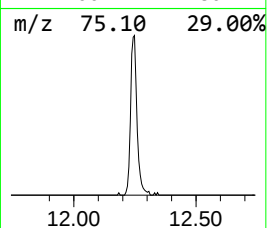
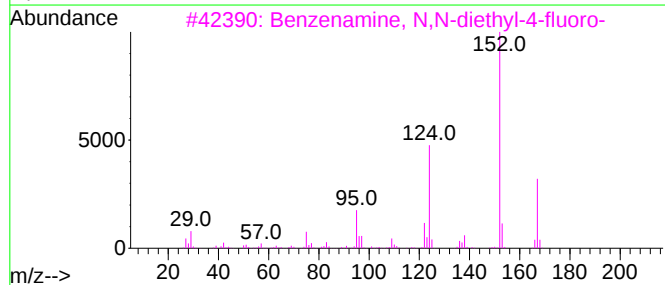
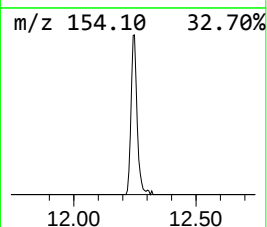
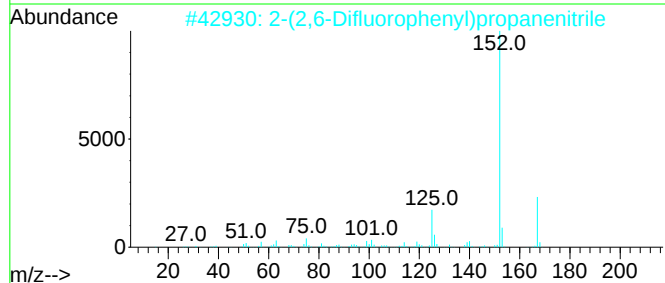
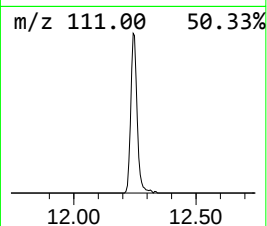
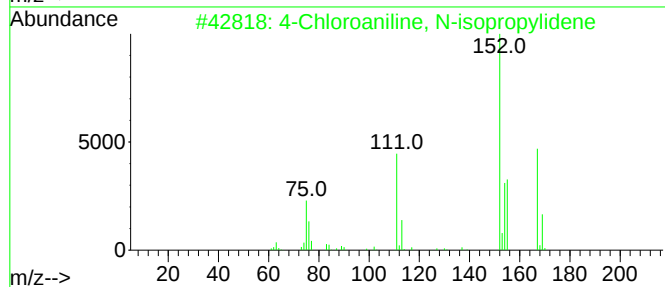
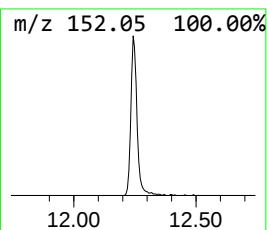
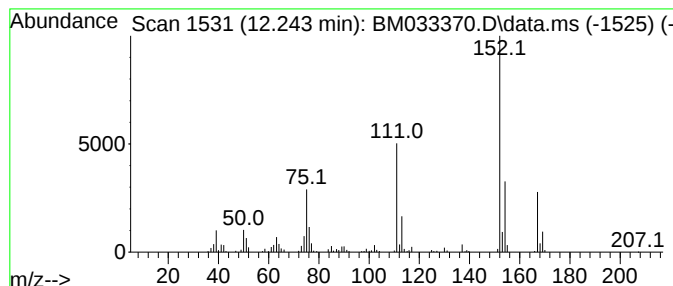
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Peak Number 1 4-Chloroaniline, N-isopropy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.243	12.82 ng/ul	275527	Naphthalene-d8	10.701	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloroaniline, N-isopropylidene	167	C9H10ClN	040938-43-0	94
2	2-(2,6-Difluorophenyl)propanenit...	167	C9H7F2N	149680-18-2	38
3	Benzenamine, N,N-diethyl-4-fluoro-	167	C10H14FN	000347-39-7	38
4	Benzonitrile, 2-amino-5-chloro-	152	C7H5ClN2	005922-60-1	38
5	5-Aminopyrimidine, N-trimethylsi...	167	C7H13N3Si	1000493-88-4	38



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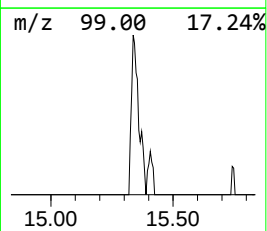
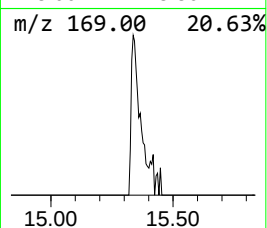
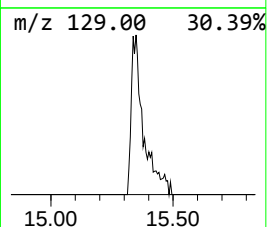
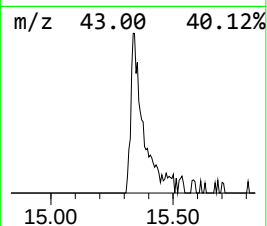
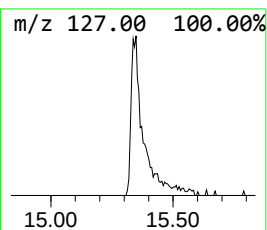
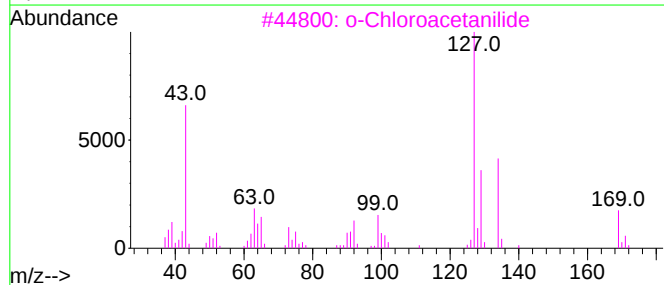
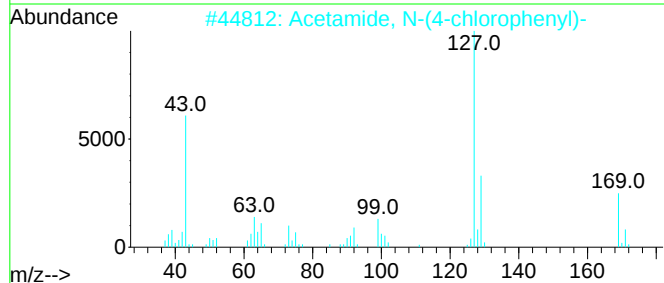
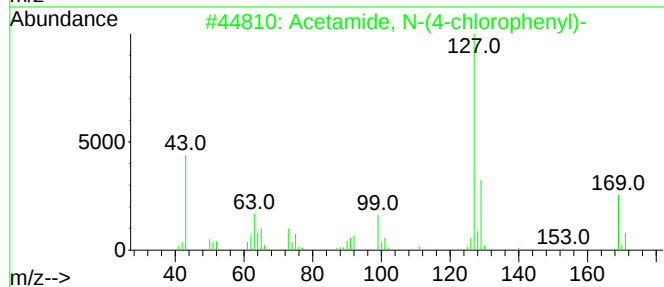
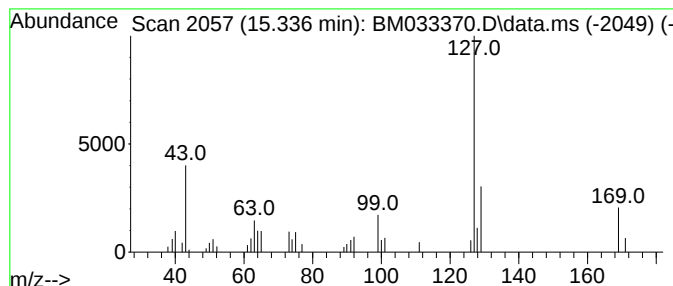
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Peak Number 2 Acetamide, N-(4-chlorophenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.336	2.63 ng/ul	70587	Acenaphthene-d10	14.531

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(4-chlorophenyl)-	169	C8H8ClNO	000539-03-7	96
2			Acetamide, N-(4-chlorophenyl)-	169	C8H8ClNO	000539-03-7	93
3			o-Chloroacetanilide	169	C8H8ClNO	000533-17-5	91
4			Acetamide, N-(4-chlorophenyl)-	169	C8H8ClNO	000539-03-7	87
5			Acetamide, N-(3-chlorophenyl)-	169	C8H8ClNO	000588-07-8	87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
4-Chloroaniline...	12.243	12.8	ng/ul	275527	2	10.701	429897	20.0
Acetamide, N-(4...	15.336	2.6	ng/ul	70587	3	14.531	537763	20.0