

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011625\
 Data File : BF141184.D
 Acq On : 16 Jan 2025 14:40
 Operator : RC/JU
 Sample : PB166033BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB166033BL

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_F\Methods\8270-BF011025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141184.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	3	9	16	rVB	104199	178886	2.55%	0.538%
2	5.428	561	567	573	rBV	2406062	3289740	46.94%	9.902%
3	6.434	731	738	743	rBV	2397929	3330079	47.51%	10.024%
4	6.810	797	802	807	rBV	707232	878224	12.53%	2.643%
5	7.375	890	898	903	rBV	2215001	3263998	46.57%	9.825%
6	8.092	1014	1020	1026	rBV	886038	1191820	17.01%	3.587%
7	9.175	1196	1204	1209	rBV	4383034	6238818	89.02%	18.779%
8	9.851	1312	1319	1324	rBV	1071826	1389894	19.83%	4.184%
9	10.639	1446	1453	1476	rBV	2002765	2637238	37.63%	7.938%
10	11.339	1566	1572	1577	rBV	1135925	1445782	20.63%	4.352%
11	11.498	1594	1599	1604	rBV	187787	232108	3.31%	0.699%
12	12.933	1836	1843	1848	rBV	5100586	7008522	100.00%	21.096%
13	13.980	2015	2021	2039	rBV	918373	1199057	17.11%	3.609%
14	15.439	2263	2269	2275	rBV	597546	937874	13.38%	2.823%

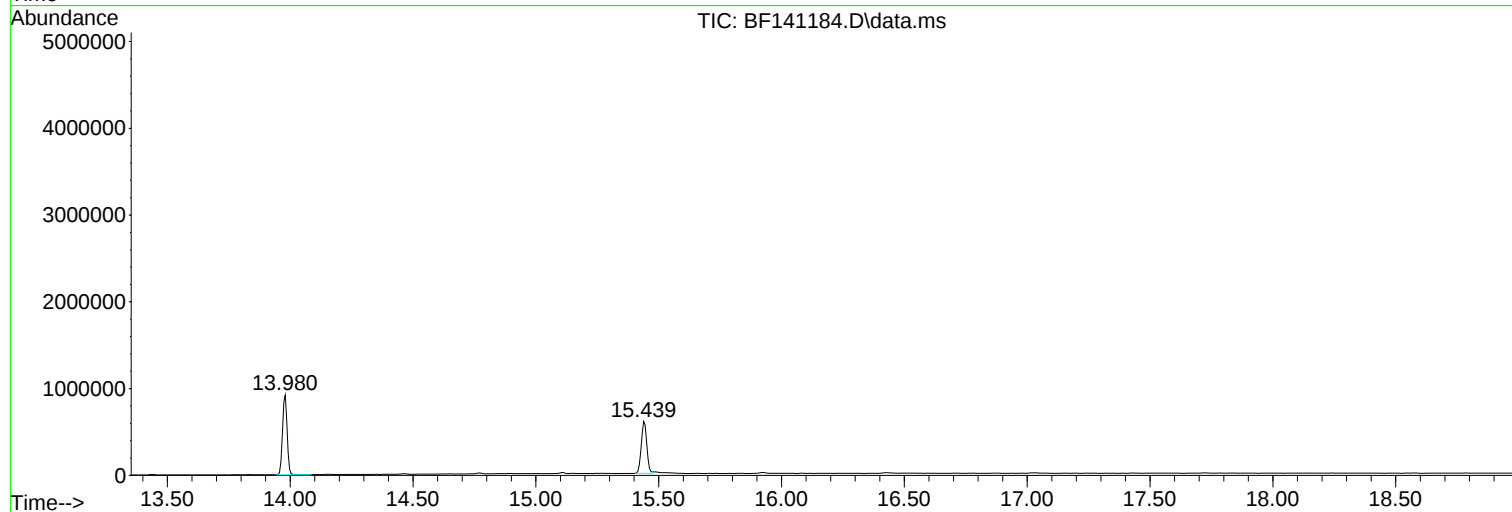
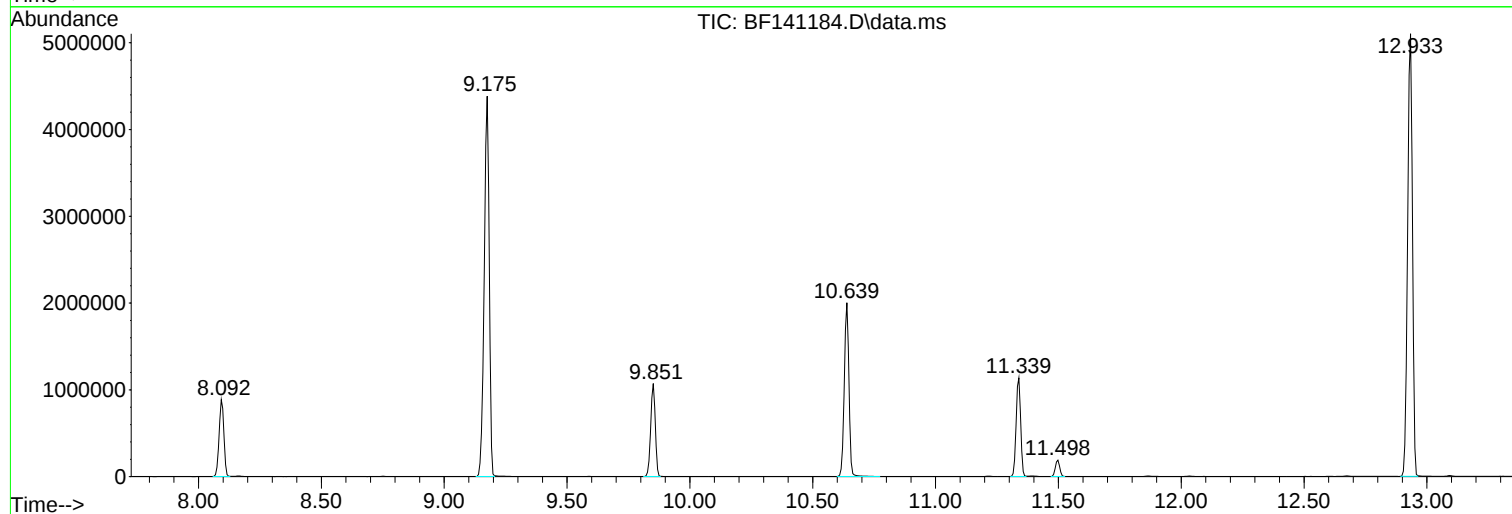
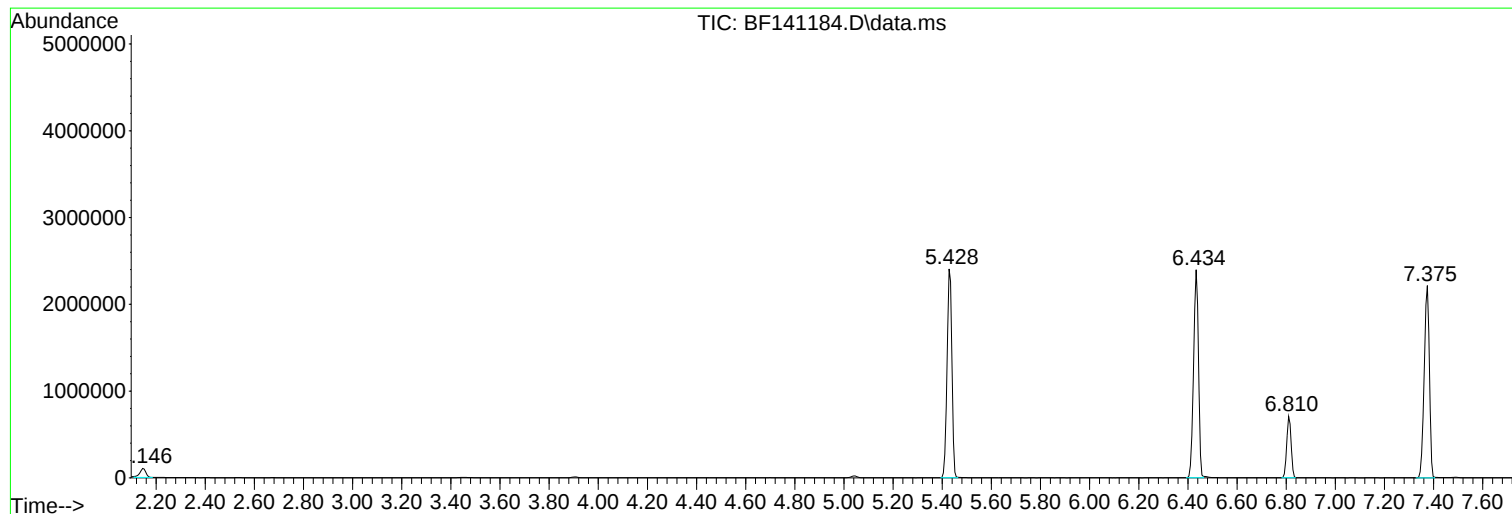
Sum of corrected areas: 33222040

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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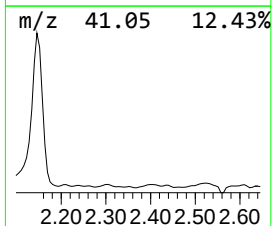
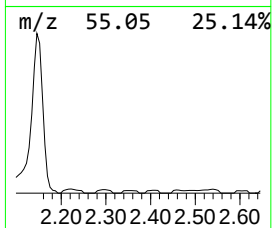
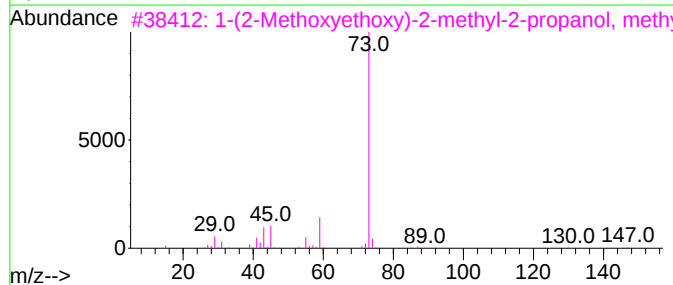
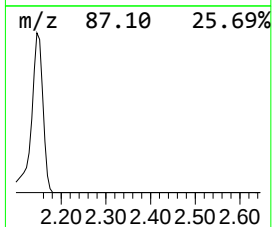
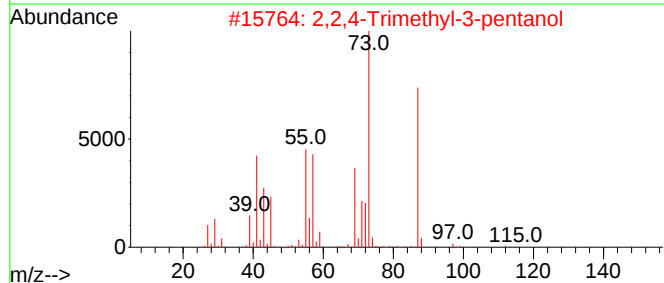
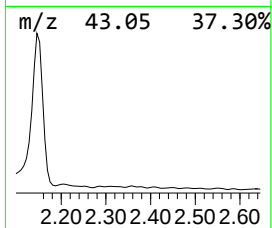
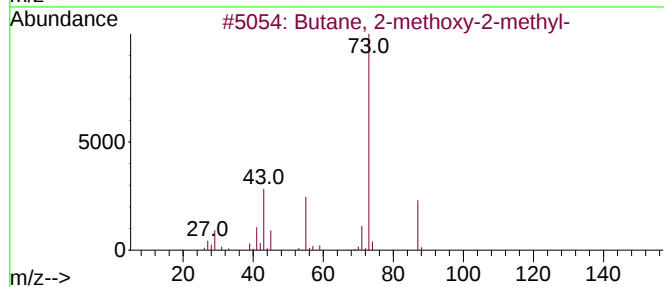
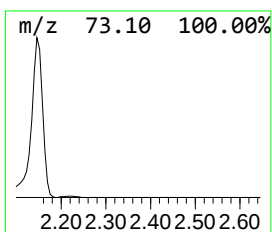
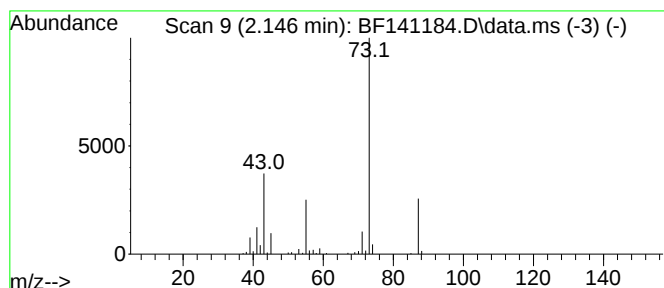
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.146	4.07 ng	178886	1,4-Dichlorobenzene-d4	6.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	25
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	23
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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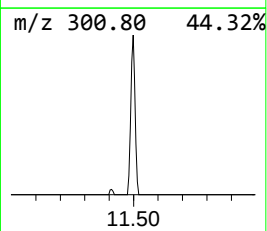
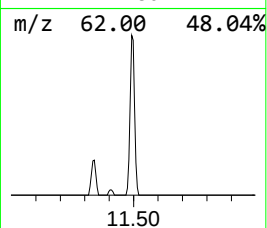
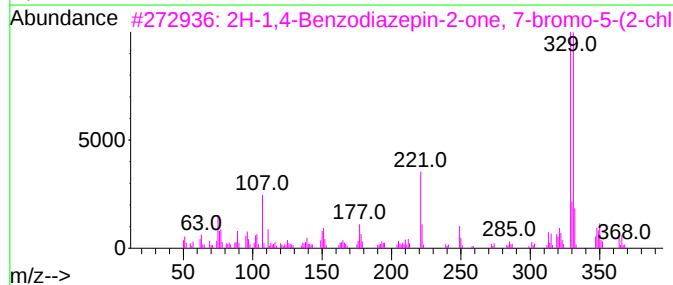
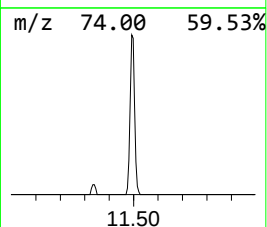
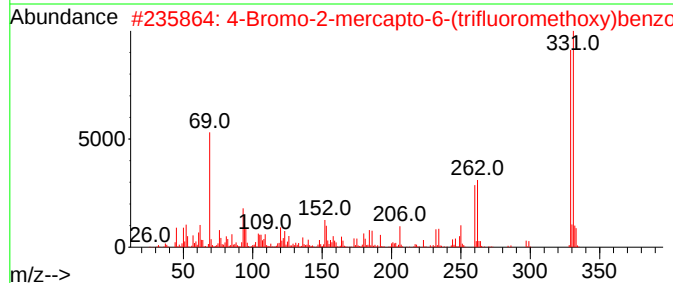
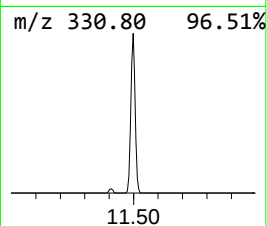
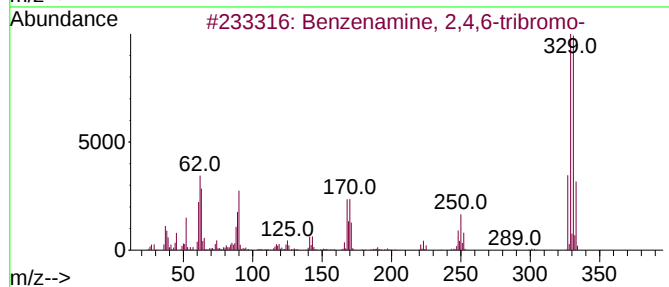
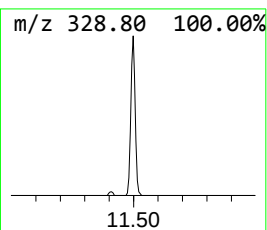
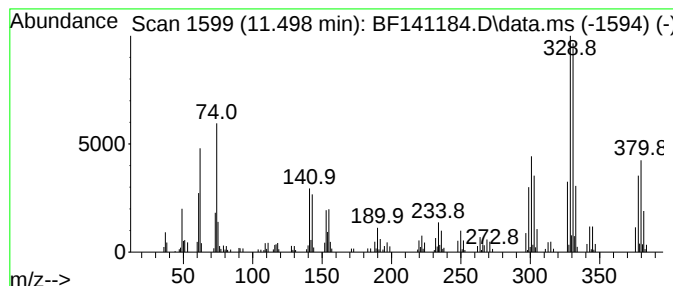
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Peak Number 2 Benzenamine, 2,4,6-tribromo- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.498	3.21 ng	232108	Phenanthrene-d10	11.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,4,6-tribromo-	327	C6H4Br3N	000147-82-0	50
2			4-Bromo-2-mercapto-6-(trifluorom...	329	C8H3BrF3NOS2	1000502-92-7	20
3			2H-1,4-Benzodiazepin-2-one, 7-br...	364	C15H10BrClN2O2	070030-09-0	12
4			Bromophos	364	C8H8BrCl2O3PS	002104-96-3	10
5			1,3,4-Oxadiazol-2-amine, 5-(2-br...	329	C11H12BrN3O4	1000350-74-8	10



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.146	4.1	ng	178886	1	6.810	878224	20.0
Benzenamine, 2,...	11.498	3.2	ng	232108	4	11.339	1445780	20.0