

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101723\
 Data File : BF135816.D
 Acq On : 17 Oct 2023 17:20
 Operator : CG\JU
 Sample : PB156288BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB156288BL

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-BF101323.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135816.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.105	50	53	58	rVB2	18177119	31774830	100.00%	56.249%
2	1.735	150	160	174	rVB	607353	830930	2.62%	1.471%
3	4.982	707	712	731	rBV	634310	1019954	3.21%	1.806%
4	5.387	777	781	803	rBV	2814709	2978496	9.37%	5.273%
5	6.399	948	953	980	rBV	2689492	3170639	9.98%	5.613%
6	6.734	1007	1010	1021	rBV	665914	601007	1.89%	1.064%
7	7.311	1104	1108	1130	rBV	2284556	2069661	6.51%	3.664%
8	8.017	1225	1228	1246	rBV	922331	819674	2.58%	1.451%
9	9.099	1407	1412	1415	rBV	3979323	3416922	10.75%	6.049%
10	9.775	1523	1527	1537	rBV	1097210	941036	2.96%	1.666%
11	10.575	1659	1663	1683	rBV	2381716	2366718	7.45%	4.190%
12	11.269	1778	1781	1790	rBV	1151092	1020993	3.21%	1.807%
13	12.875	2048	2054	2057	rBV	3703888	3826442	12.04%	6.774%
14	13.940	2231	2235	2246	rBV	849341	879160	2.77%	1.556%
15	15.410	2481	2485	2496	rBV	521766	773492	2.43%	1.369%

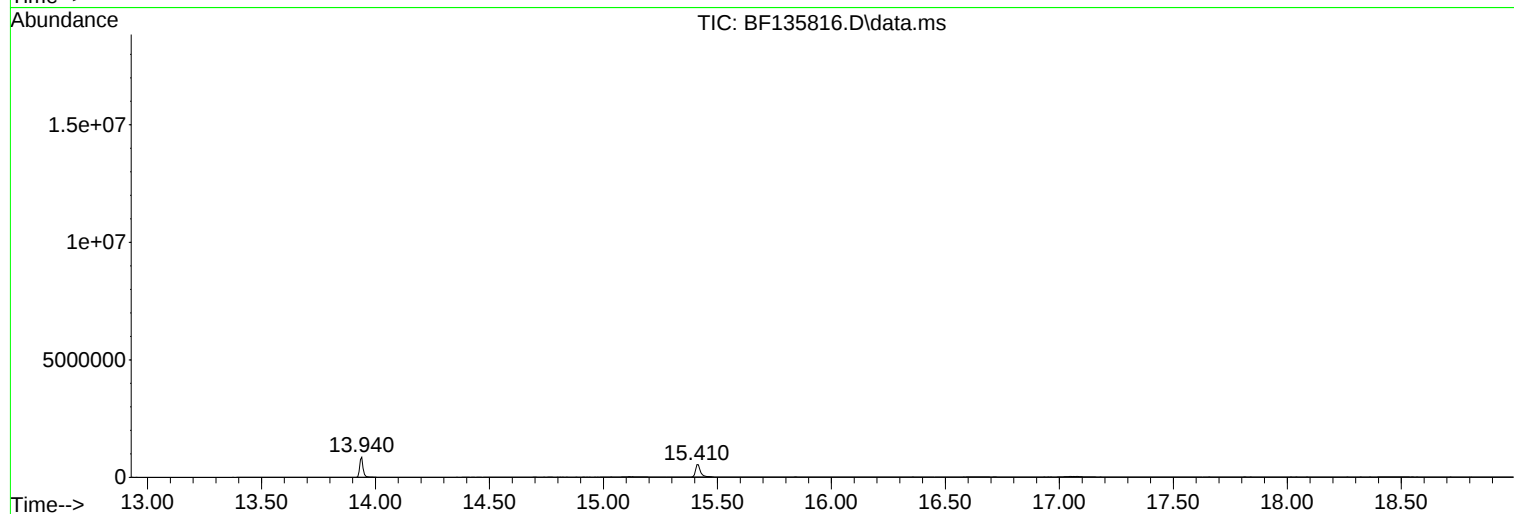
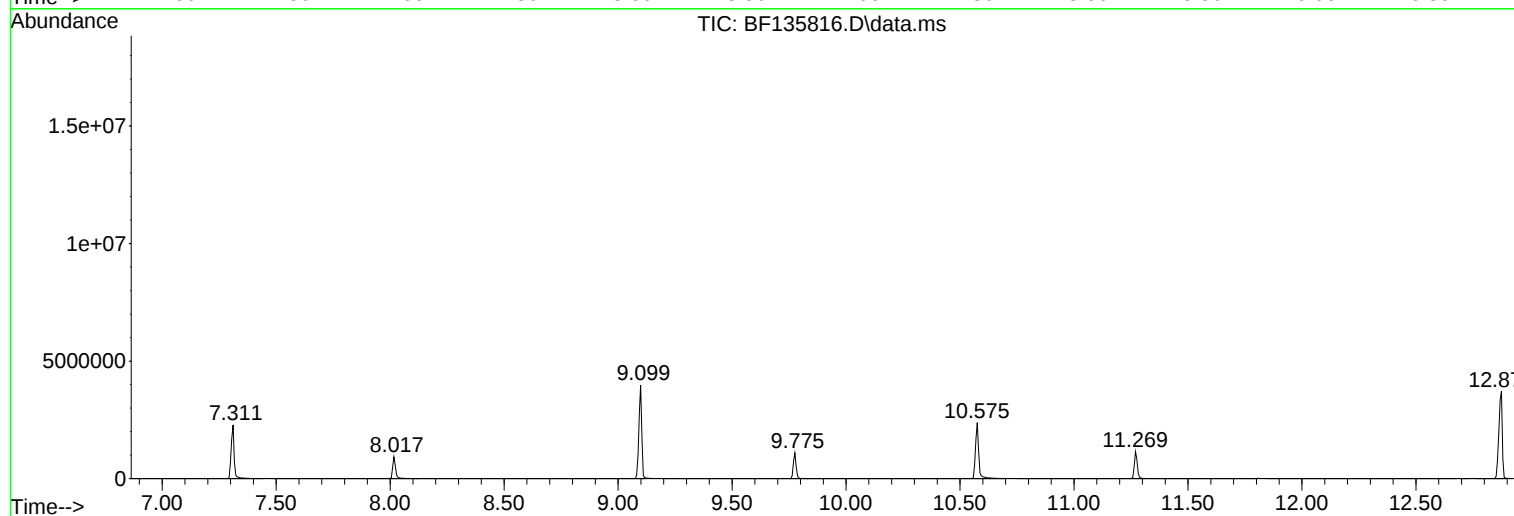
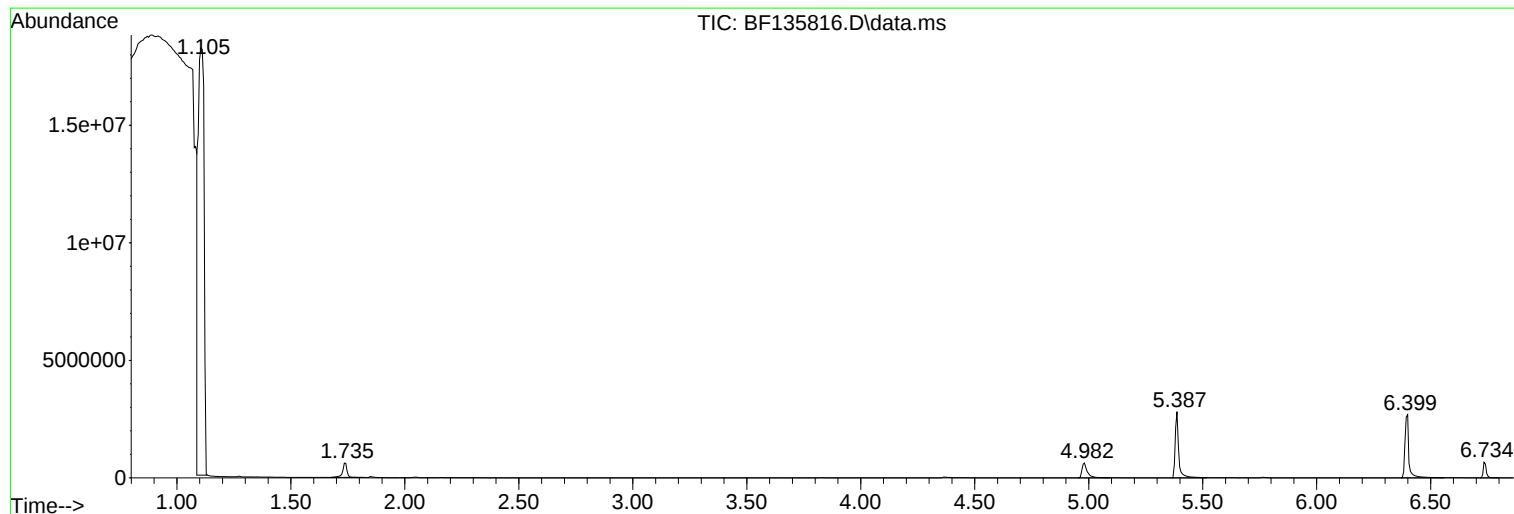
Sum of corrected areas: 56489954

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.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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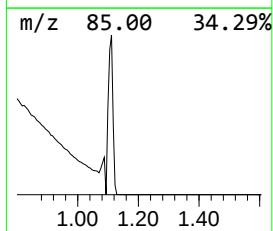
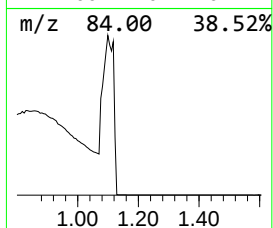
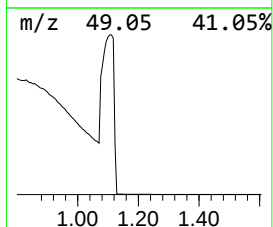
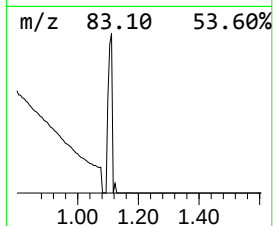
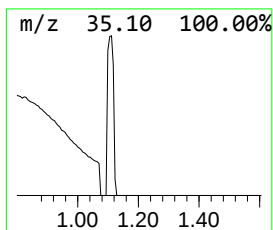
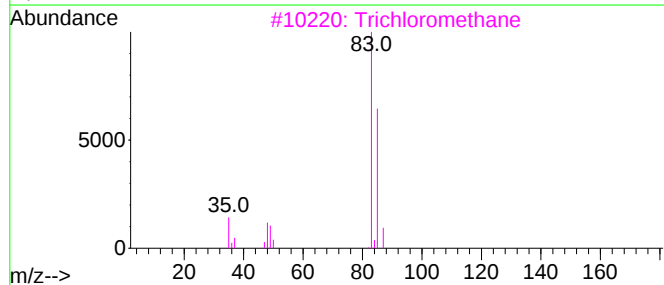
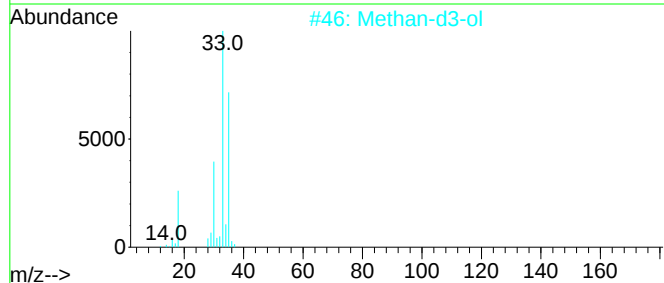
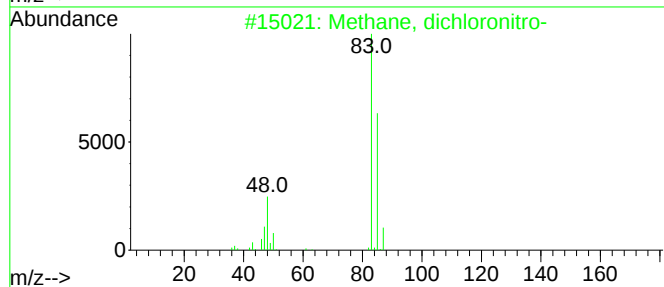
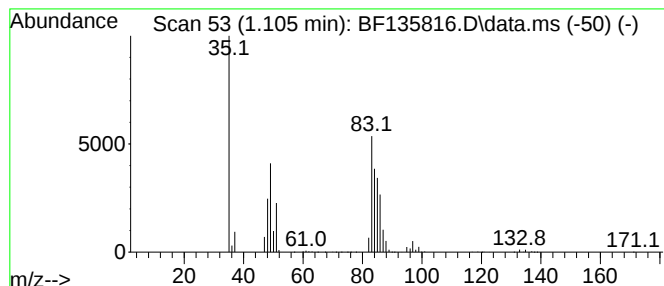
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.105 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.105	1057.39 ng	31774800	1,4-Dichlorobenzene-d4	6.734

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methane, dichloronitro-	129	CHCl2NO2	007119-89-3	9
2			Methan-d3-ol	35	CHD3O	001849-29-2	4
3			Trichloromethane	118	CHCl3	000067-66-3	4
4			Thionyl chloride	118	Cl2OS	007719-09-7	4
5			Ethanol, 2,2-dichloro-	114	C2H4Cl2O	000598-38-9	2



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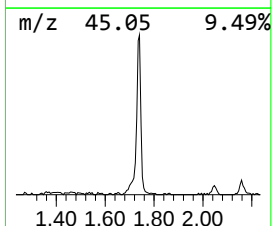
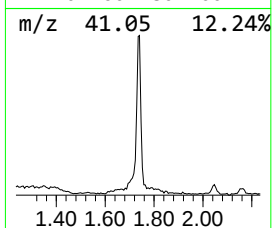
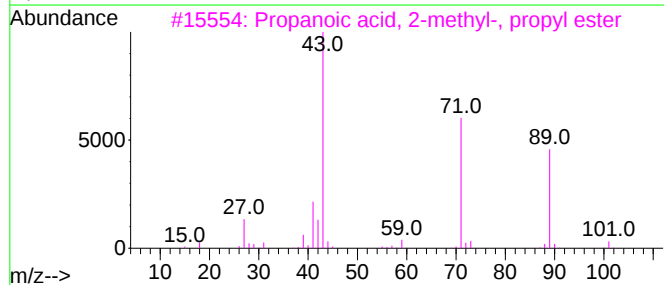
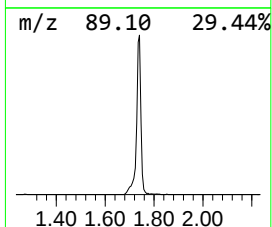
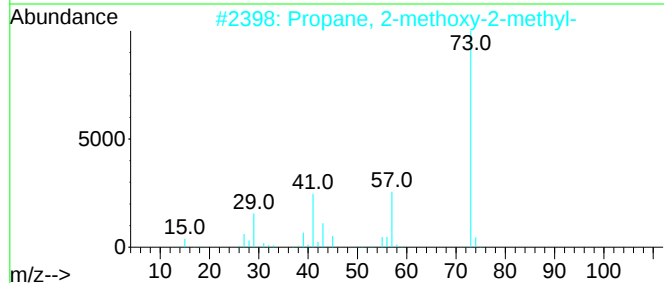
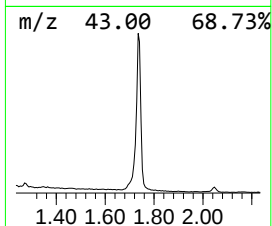
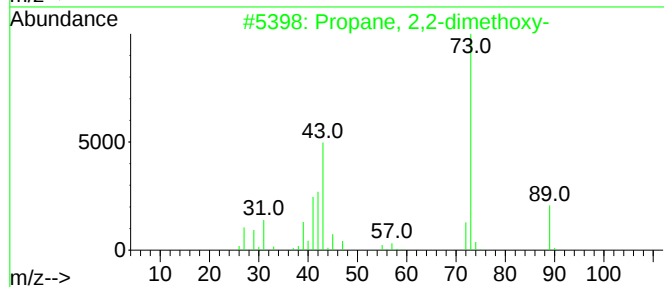
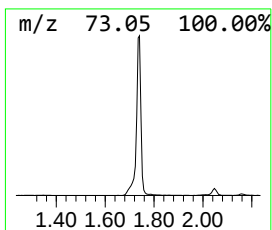
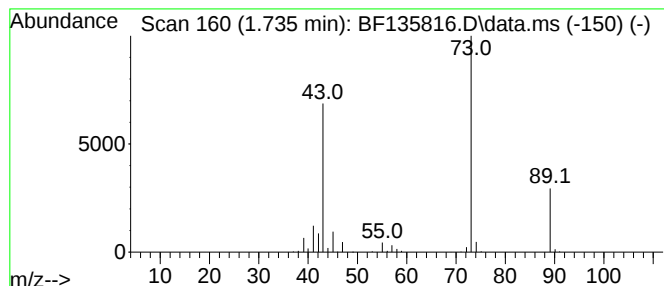
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Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.735	27.65 ng	830930	1,4-Dichlorobenzene-d4	6.734

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	23
3			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	23
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	9



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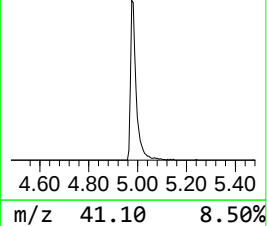
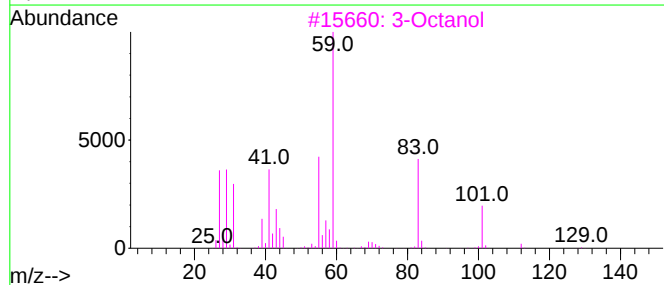
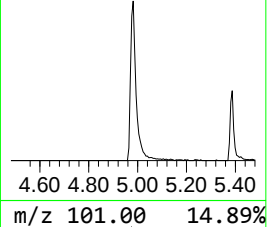
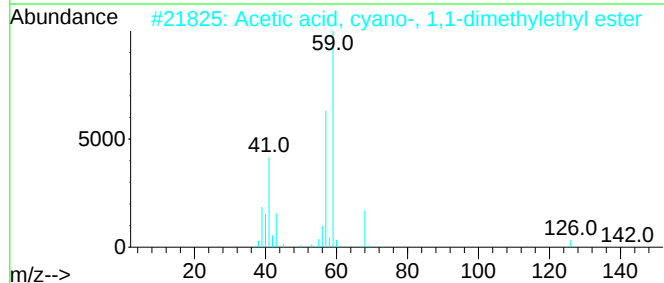
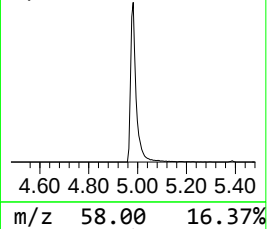
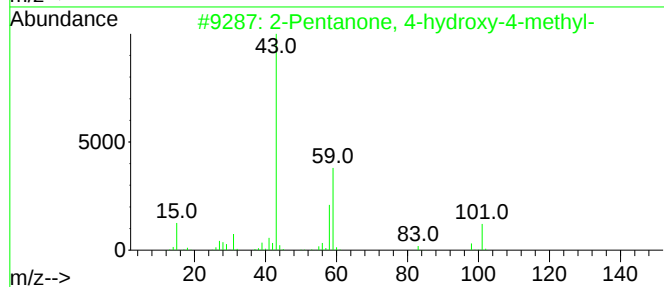
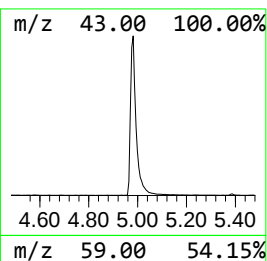
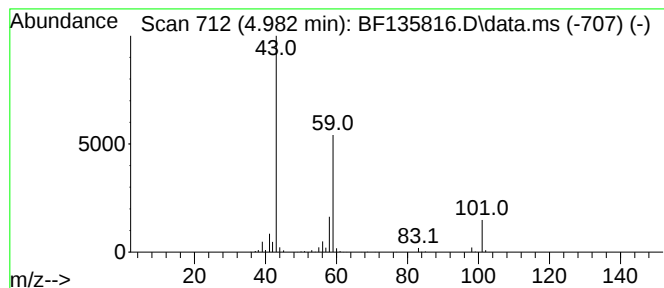
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.982	33.94 ng	1019950	1,4-Dichlorobenzene-d4			6.734
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, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	37
3	3-Octanol		130	C8H18O	000589-98-0	33
4	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	33
5	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	28



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.105	1.105	1057.4	ng	31774800	1	6.734	601007	20.0
Propane, 2,2-di...	1.735	27.6	ng	830930	1	6.734	601007	20.0
2-Pentanone, 4-...	4.982	33.9	ng	1019950	1	6.734	601007	20.0