Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
Data File : BP011093.D
Acq On : 25 Jul 2022 15:35
Operator : CG/JU
Sample : N3801-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 25 23:21:16 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 19 03:25:54 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
GBHW8

Manual IntegrationsAPPROVED

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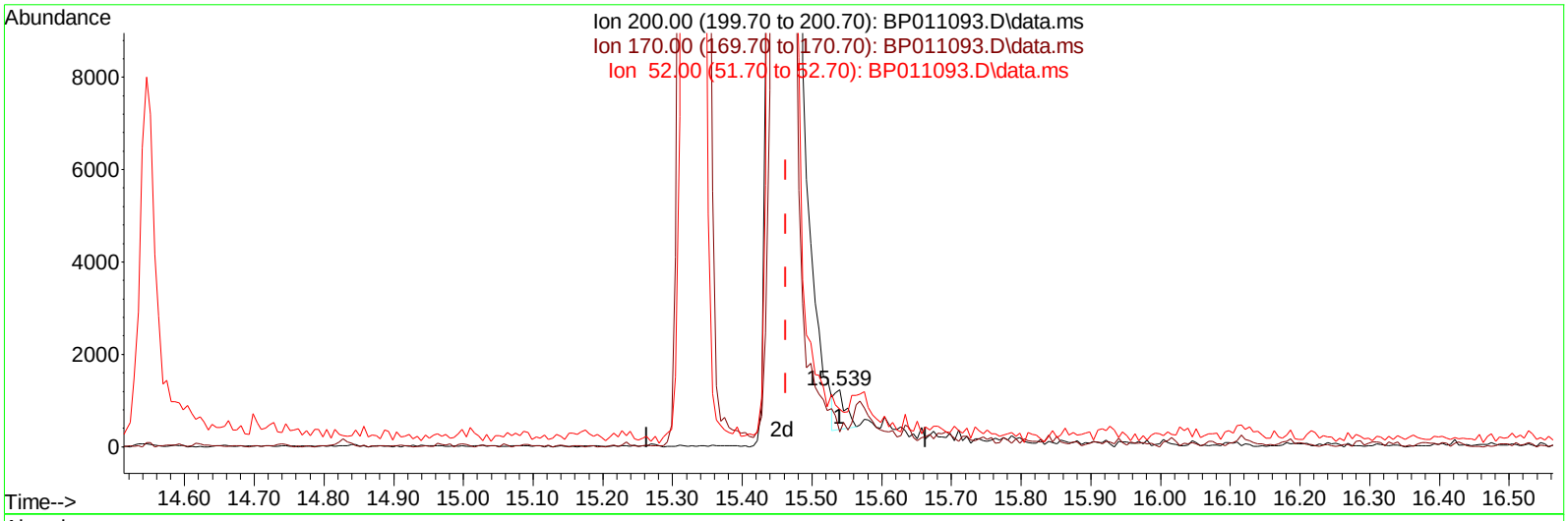
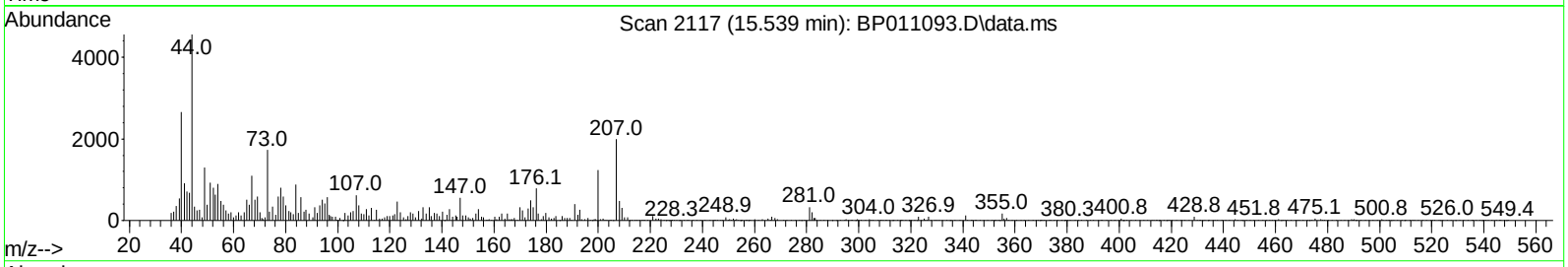
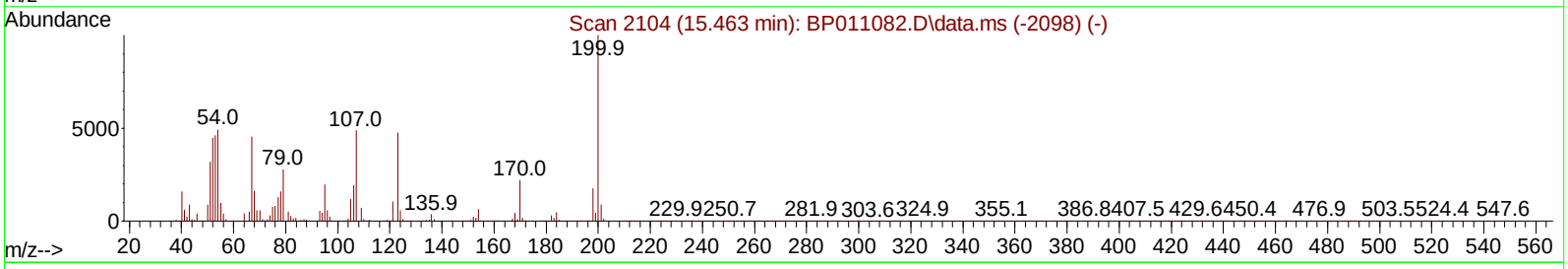
Supervised By :mohammad ahmed 07/26/2022







TIC: BP011093.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.539min (+ 0.076) 0.08 ng/ul

response 951

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	26.26
52.00	57.40	65.61
0.00	0.00	0.00

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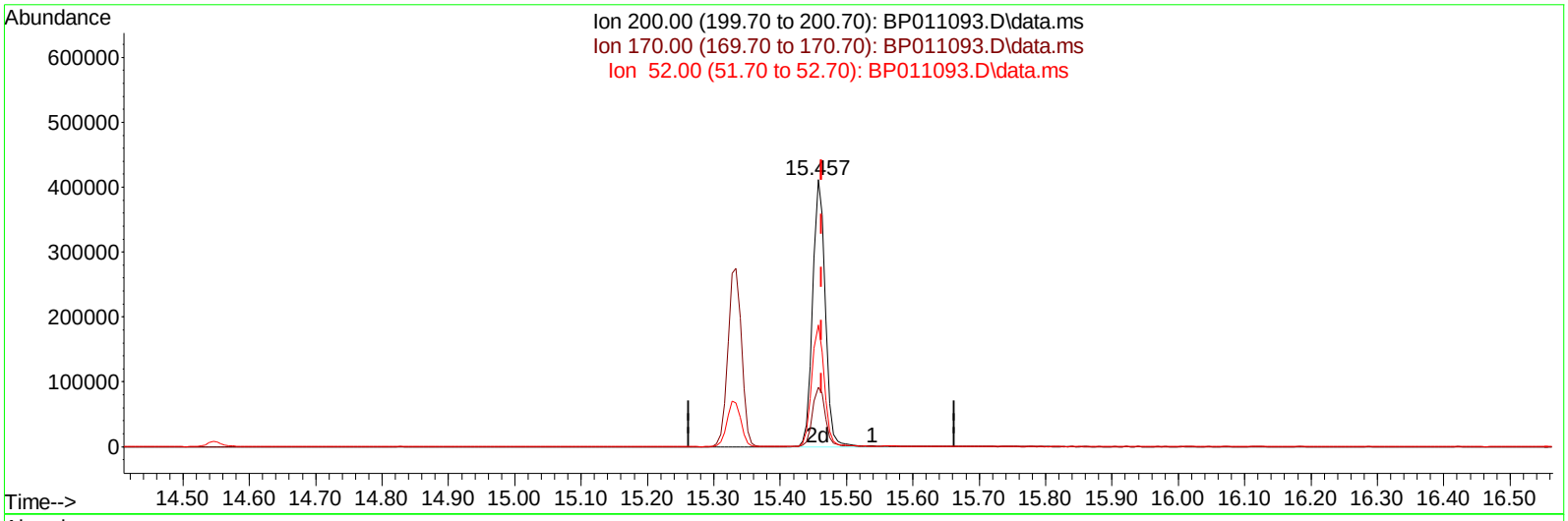

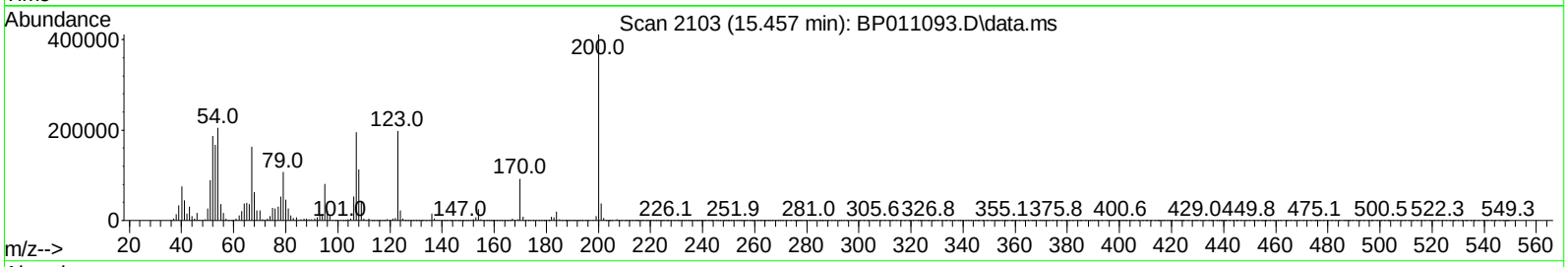

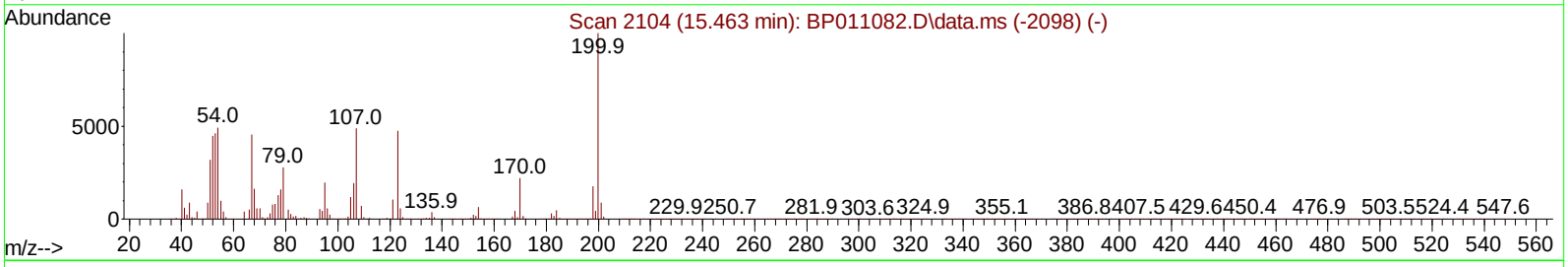

TIC: BP011093.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.457min (-0.006) 48.03 ng/ul m

response 544254

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	22.40
52.00	57.40	45.57#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.669	152	378330	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.457	136	1796160	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.328	164	1131891	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.098	188	2289878	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.216	240	1836677	20.000	ng/ul	#-0.01
88) Perylene-d12	23.557	264	1462633	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	55715	5.309	ng/uL	-0.02
4) Pyridine-d5	3.575	84	239094	8.815	ng/ul	-0.02
7) Phenol-d5	6.846	99	269932	8.692	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	907598	45.448	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.199	132	829986	33.075	ng/ul	-0.01
15) 4-Methylphenol-d8	8.381	113	569933	22.924	ng/ul	-0.02
21) Nitrobenzene-d5	8.828	128	598563	45.728	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.552	143	605083	46.717	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	916192	38.680	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.610	131	1367890	35.030	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	3967183	51.282	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	4605447	50.456	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	91821	5.865	ng/ul	-0.01
60) Fluorene-d10	15.334	176	3242731	49.314	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	544254m	48.032	ng/ul	0.00
73) Anthracene-d10	17.198	188	4981440	51.171	ng/ul	0.00
81) Pyrene-d10	19.451	212	5425203	56.242	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.404	264	3726493	53.047	ng/ul	-0.02
Target Compounds						Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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