Quantitation Report (Qedit)

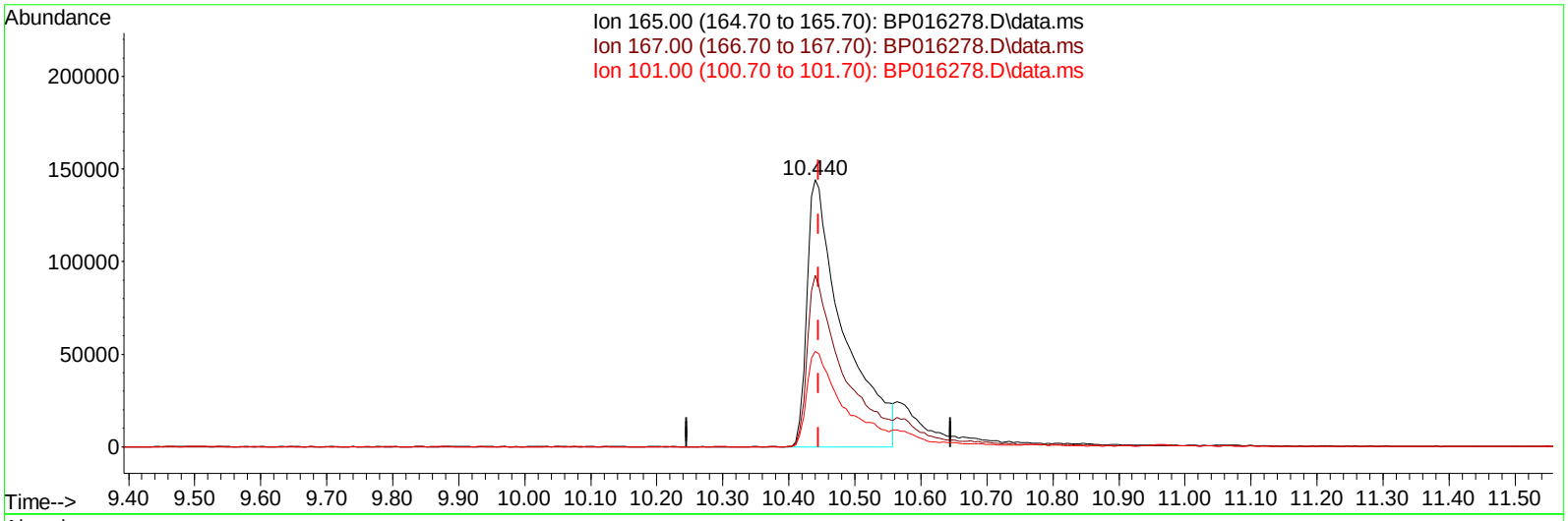
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071723\
Data File : BP016278.D
Acq On : 18 Jul 2023 05:43
Operator : MA/JU
Sample : PB154017BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

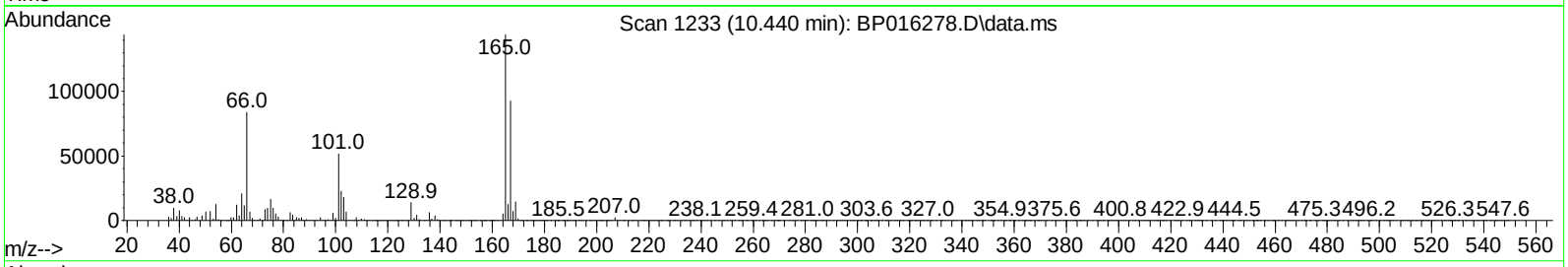
Quant Time: Jul 18 06:56:17 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 14 23:53:26 2023
Response via : Initial Calibration

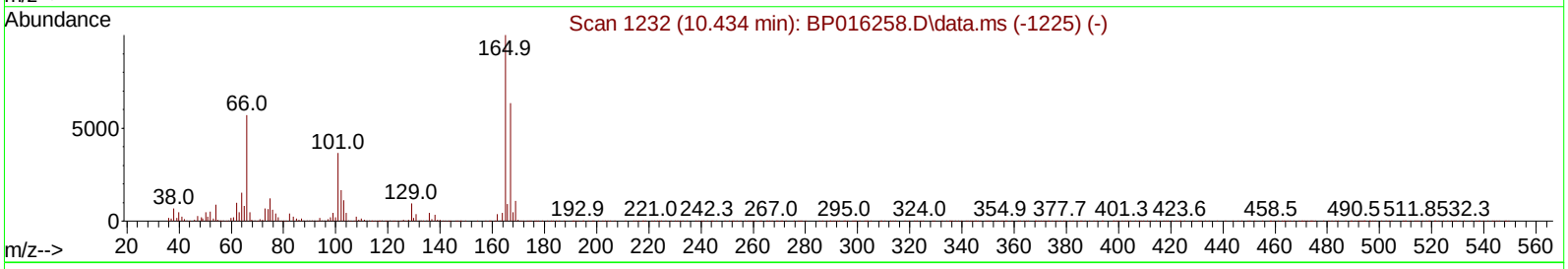
Instrument :
BNA_P
ClientSampleId :
SBLK017

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/19/2023
Supervised By :mohammad ahmed 07/19/2023







TIC: BP016278.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.440min (-0.006) 33.94 ng/ul

response 550383

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.20	64.33
101.00	38.50	35.80
0.00	0.00	0.00

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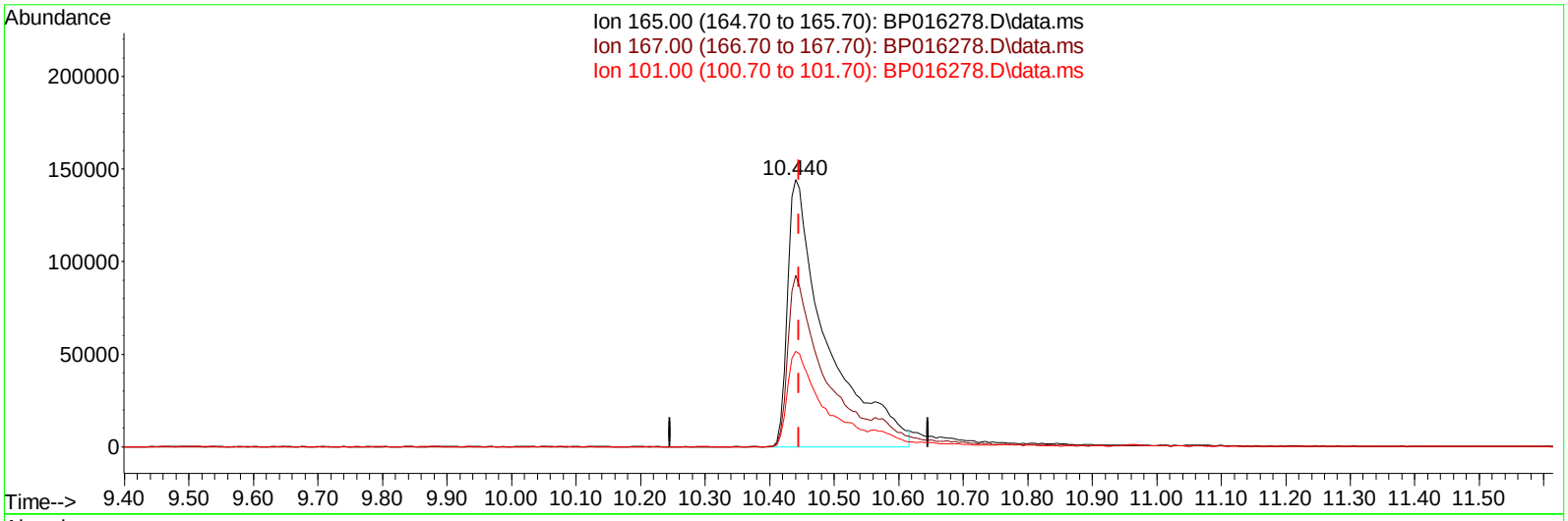

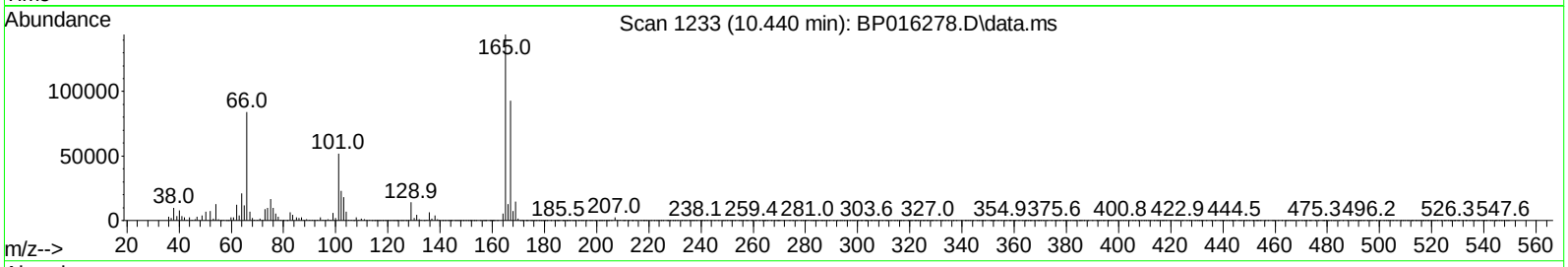

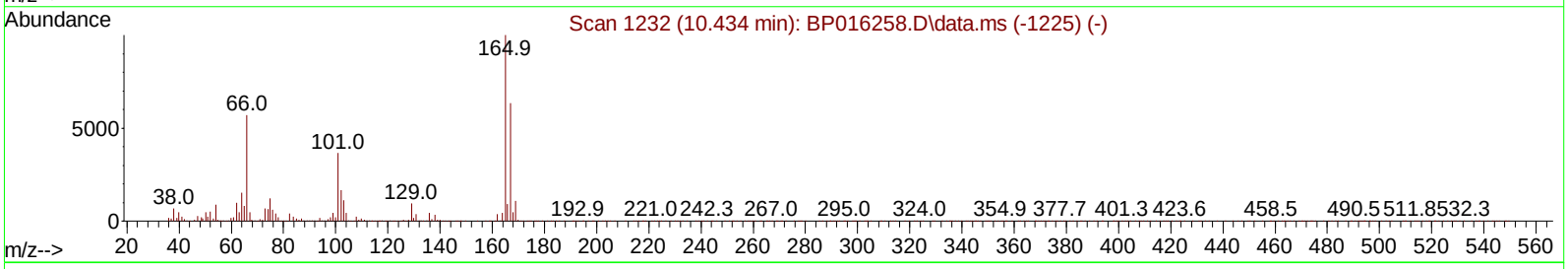

TIC: BP016278.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.440min (-0.006) 37.49 ng/ul m

response 607942

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	65.20	64.33
101.00	38.50	35.80
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Acq On : 18 Jul 2023 05:43
Operator : MA/JU
Sample : PB154017BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Quant Time: Jul 18 06:56:58 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	255408	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	1190832	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	761278	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	1750951	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	1777052	20.000	ng/ul	0.01
88) Perylene-d12	23.963	264	1940336	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	46907	6.499	ng/uL	0.00
4) Pyridine-d5	3.728	84	623339	34.628	ng/ul	-0.01
7) Phenol-d5	7.140	99	862493	37.386	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	571985	36.667	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.481	132	649700	39.013	ng/ul	-0.01
15) 4-Methylphenol-d8	8.693	113	716518	38.555	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	331600	40.363	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	372964	39.832	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	607942m	37.493	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	869411	33.452	ng/ul	0.00
46) Dimethylphthalate-d6	14.028	166	2194585	37.791	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	2439616	38.334	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	284867	24.037	ng/ul	-0.02
60) Fluorene-d10	15.610	176	1897739	39.048	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	273730	27.081	ng/ul	0.00
73) Anthracene-d10	17.475	188	3151638	41.149	ng/ul	0.00
81) Pyrene-d10	19.710	212	3891958	42.207	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	3823216	41.030	ng/ul	0.00
Target Compounds						Qvalue

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

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