Quantitation Report (Qedit)

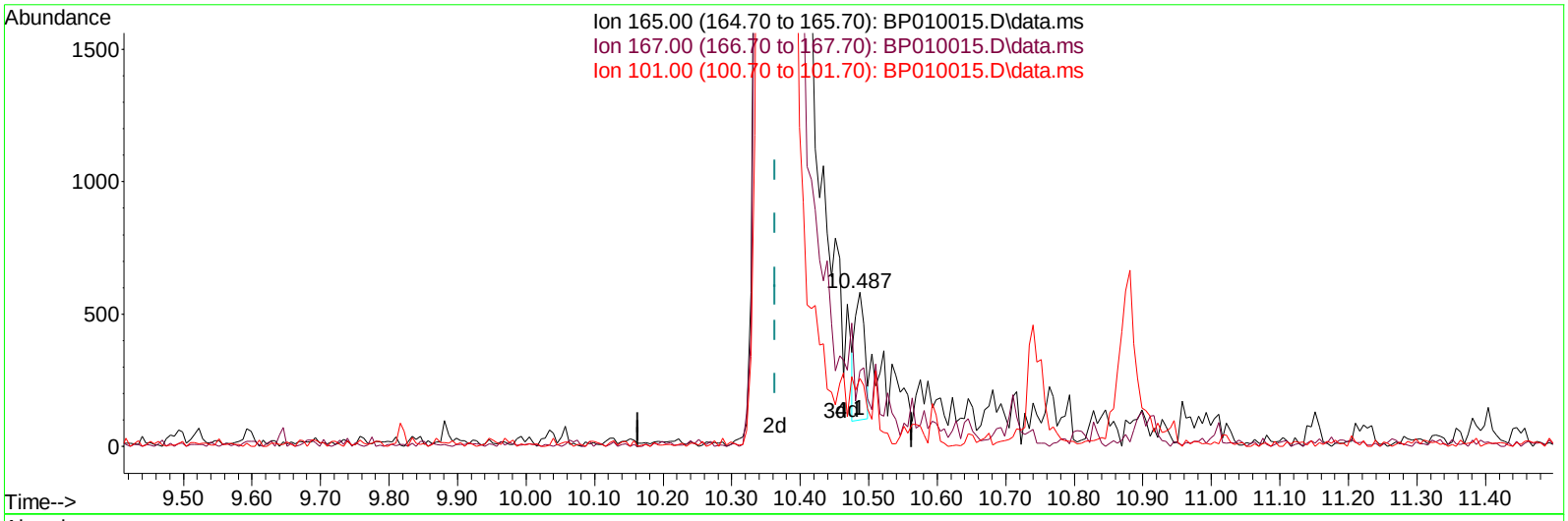
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010015.D
Acq On : 21 Apr 2022 15:37
Operator : CG/JU
Sample : PB144194BL
Misc :
ALS Vial : 10 Sample Multiplier: 1

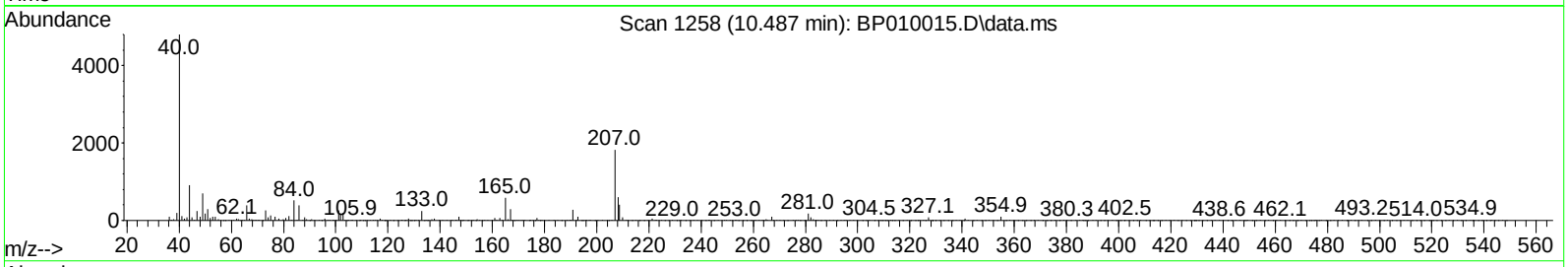
Quant Time: Apr 21 19:05:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

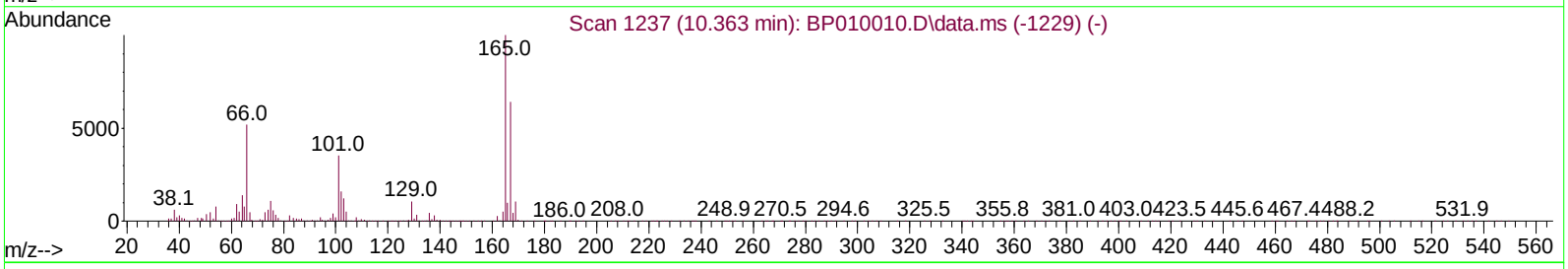
Instrument :
BNA_P
ClientSampleId :
SBLK194

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :mohammad ahmed 04/26/2022







TIC: BP010015.D\data.ms

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

10.487min (+ 0.124) 0.05 ng/ul

response 481

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	49.31
101.00	49.80	44.33
0.00	0.00	0.00

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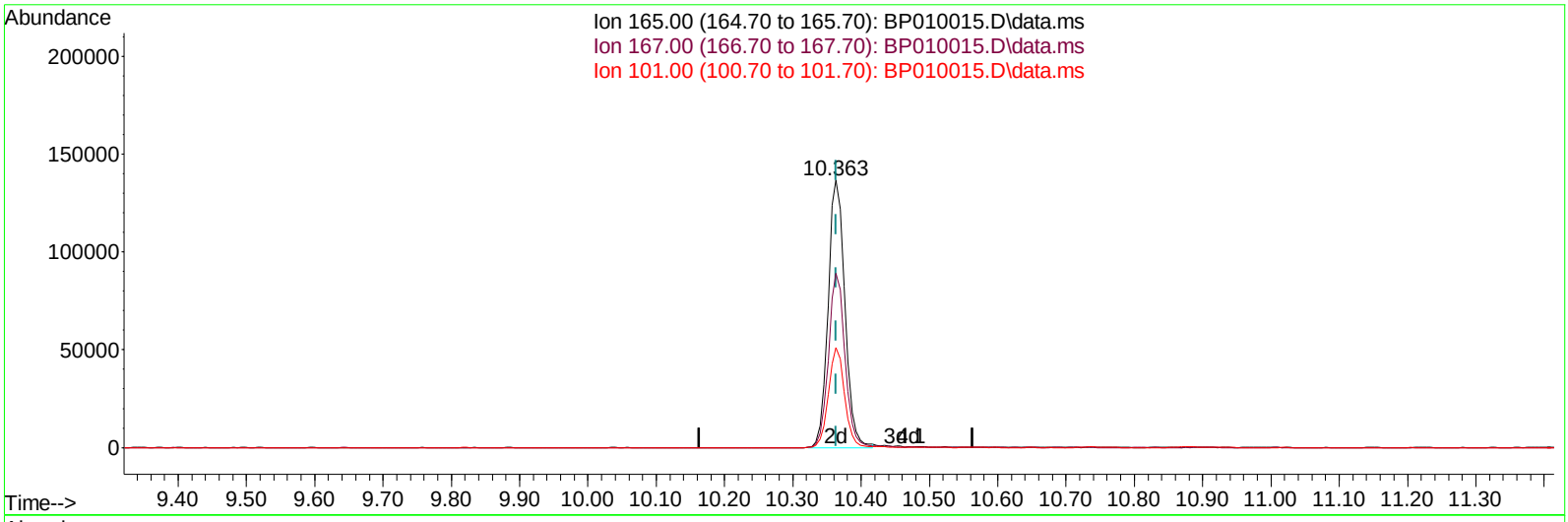

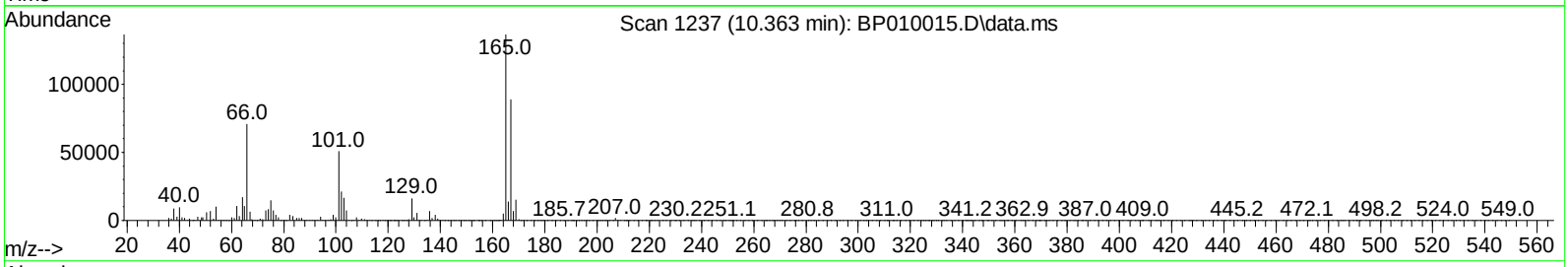

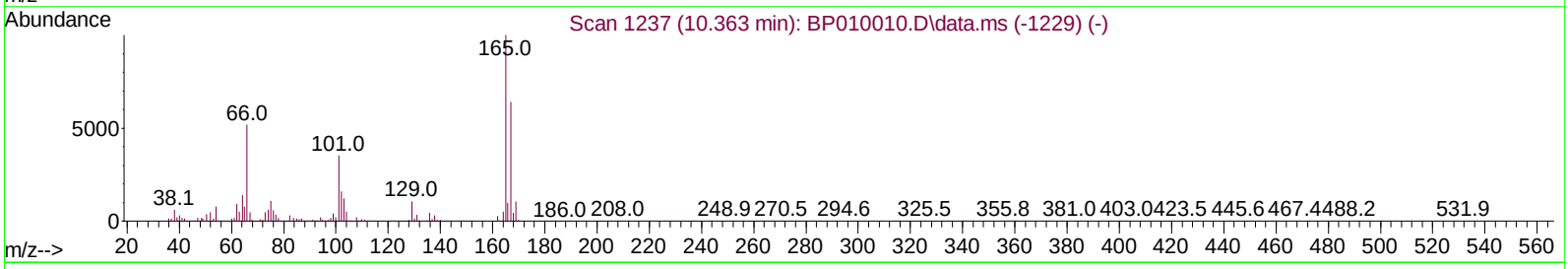

TIC: BP010015.D\data.ms

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

10.363min (+ 0.000) 23.38 ng/ul m

response 234475

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	65.10
101.00	49.80	37.28#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Supervised By :mohammad ahmed 04/26/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	129057	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	593126	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	452246	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1014985	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	1010561	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	933523	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	16983	5.738	ng/uL	0.00
4) Pyridine-d5	3.787	84	216292	25.978	ng/ul	0.00
7) Phenol-d5	7.099	99	277174	23.611	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	191850	27.236	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	227710	26.192	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	244347	25.123	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	122681	26.006	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	134655	26.136	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	234475m	23.378	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	363712	25.291	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	965786	27.453	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1052063	24.710	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	144224	22.269	ng/ul	0.00
60) Fluorene-d10	15.563	176	814907	26.929	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	135662	21.138	ng/ul	0.00
73) Anthracene-d10	17.422	188	1327188	27.028	ng/ul	0.00
81) Pyrene-d10	19.651	212	1610956	28.031	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1342459	27.430	ng/ul	0.00
Target Compounds						Qvalue

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

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