Quantitation Report (Qedit)

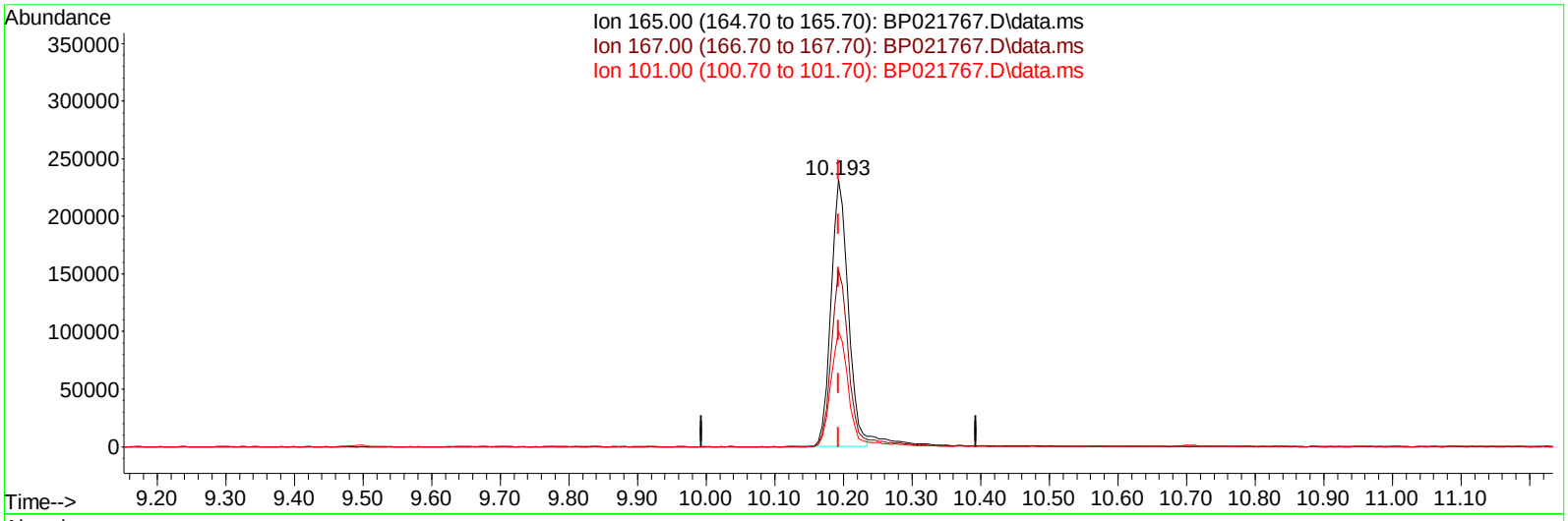
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021767.D
Acq On : 31 Aug 2024 00:59
Operator : RC/JU
Sample : P3733-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

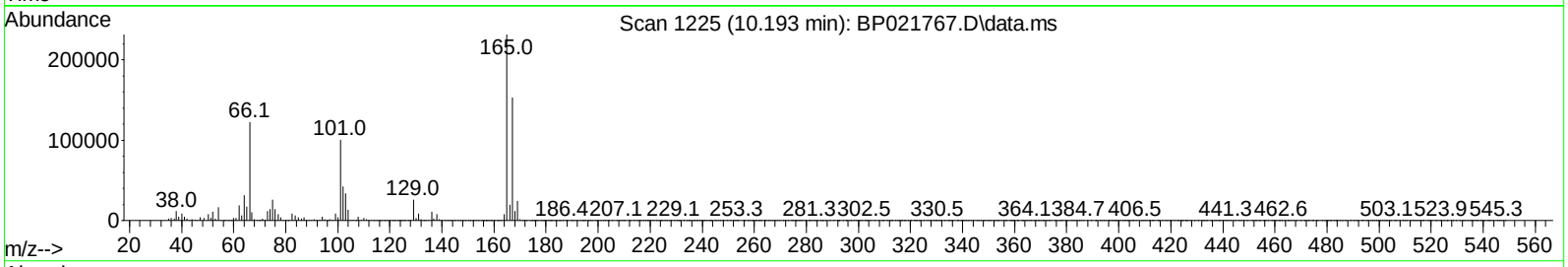
Quant Time: Aug 31 01:25:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 29 23:56:59 2024
Response via : Initial Calibration

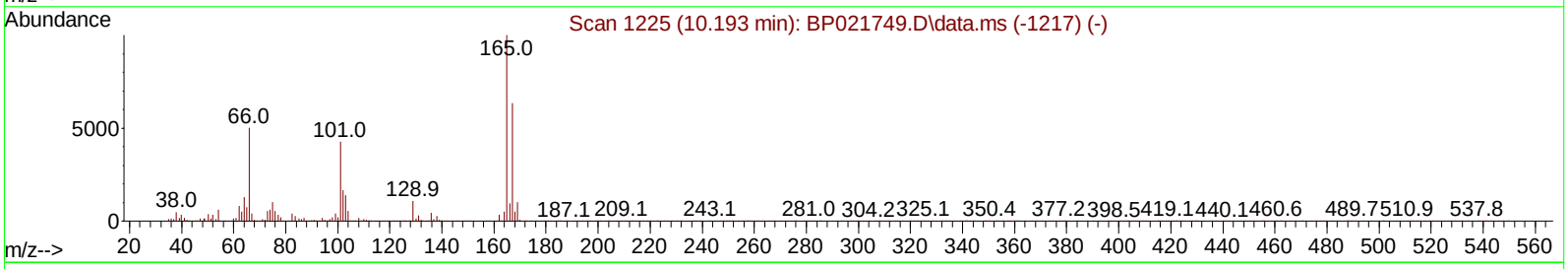
Instrument :
BNA_P
ClientSampleId :
A4499

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/02/2024
Supervised By :mohammad ahmed 09/03/2024







TIC: BP021767.D\data.ms

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

10.193min (-0.000) 22.43 ng/ul

response 403522

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	66.08
101.00	41.40	43.31
0.00	0.00	0.00

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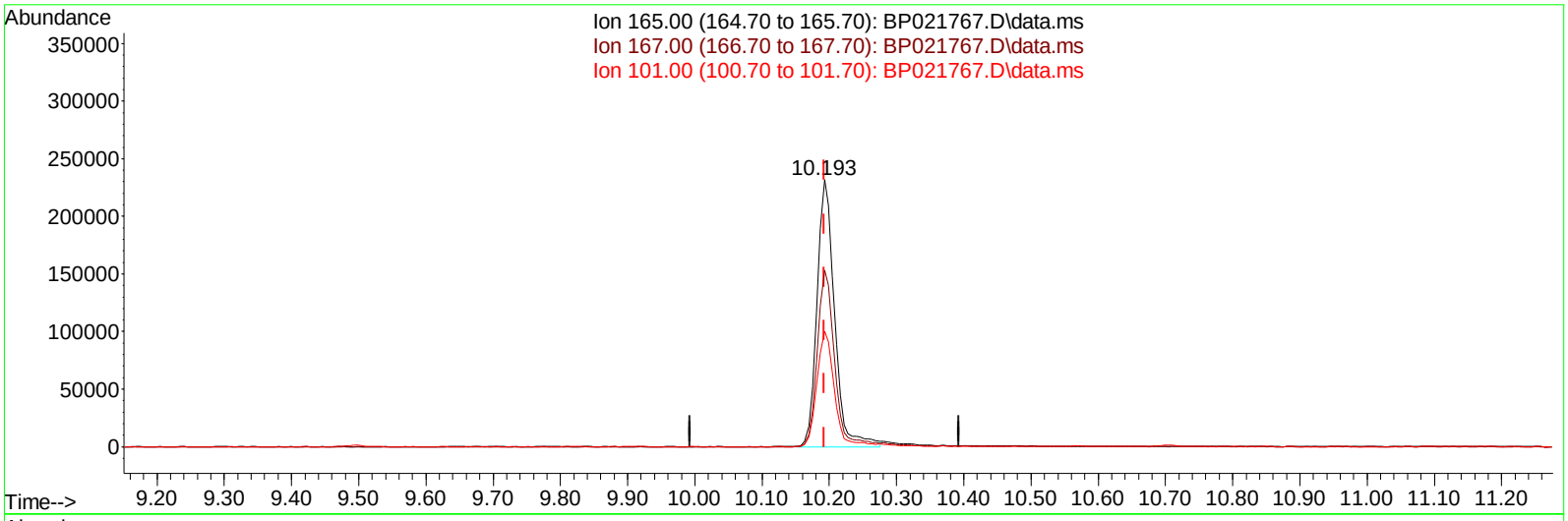

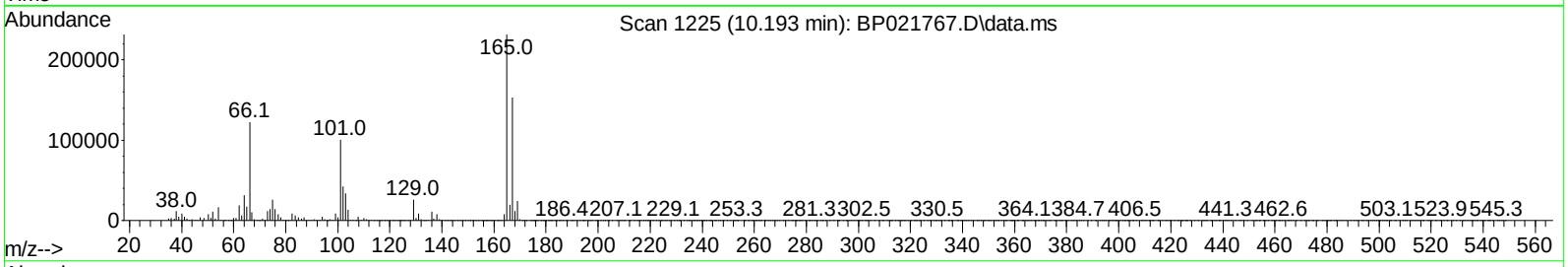

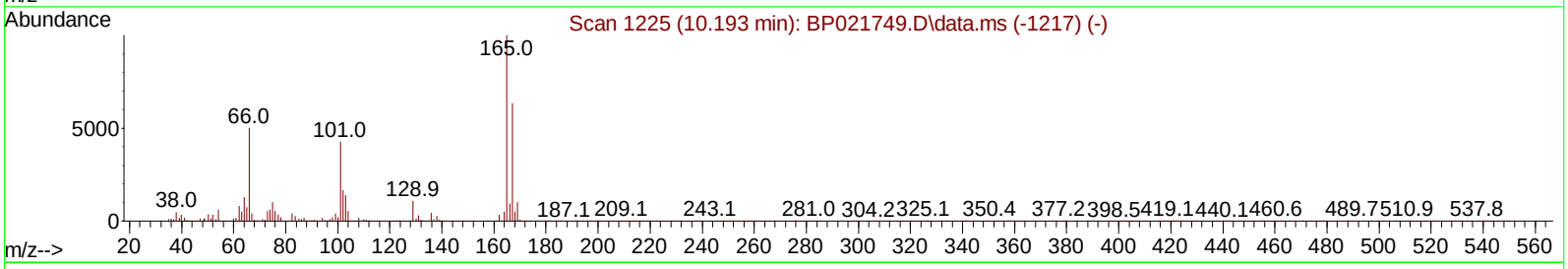

TIC: BP021767.D\data.ms

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

10.193min (-0.000) 23.40 ng/ul m

response 420889

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	66.08
101.00	41.40	43.31
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021767.D
Acq On : 31 Aug 2024 00:59
Operator : RC/JU
Sample : P3733-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 31 01:26:23 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 29 23:56:59 2024
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
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09/03/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	269666	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	1117662	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	753186	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1673810	20.000	ng/ul	0.00
79) Chrysene-d12	21.674	240	1732066	20.000	ng/ul	0.00
88) Perylene-d12	25.080	264	1959952	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	34333	4.227	ng/uL	0.00
4) Pyridine-d5	3.687	84	520559	22.157	ng/ul	0.00
7) Phenol-d5	6.958	99	687691	23.896	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	385260	24.599	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	465118	25.033	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	541476	24.349	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	232518	25.858	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	242068	26.521	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	420889m	23.400	ng/ul	0.00
31) 4-Chloroaniline-d4	10.699	131	669096	25.569	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1655423	28.724	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	1758336	27.175	ng/ul	0.00
54) 4-Nitrophenol-d4	14.634	143	269623	24.590	ng/ul	0.00
60) Fluorene-d10	15.434	176	1380711	28.496	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.551	200	190555	20.699	ng/ul	0.00
73) Anthracene-d10	17.322	188	2266720	28.535	ng/ul	-0.02
81) Pyrene-d10	19.692	212	2824837	28.530	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.845	264	3026265	28.956	ng/ul	0.00

Target Compounds Qvalue

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

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