

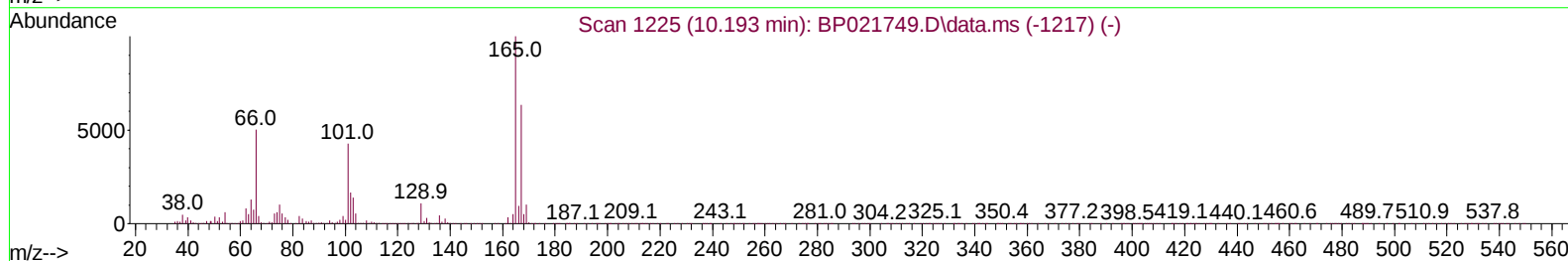
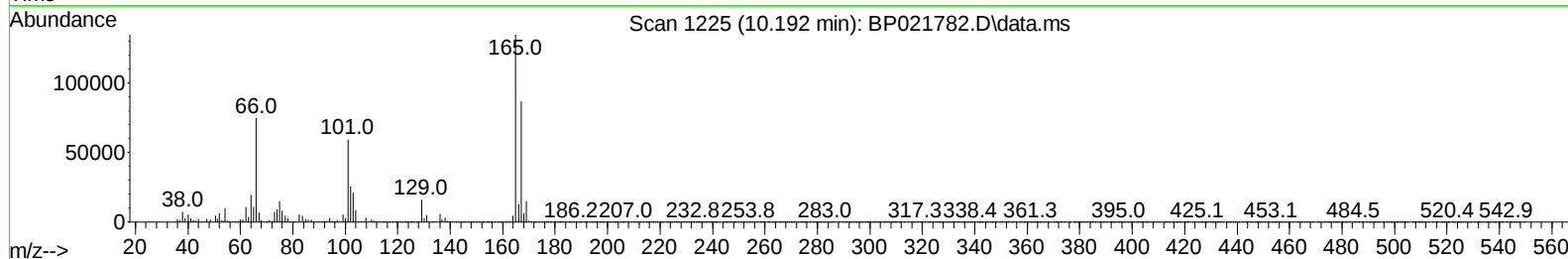
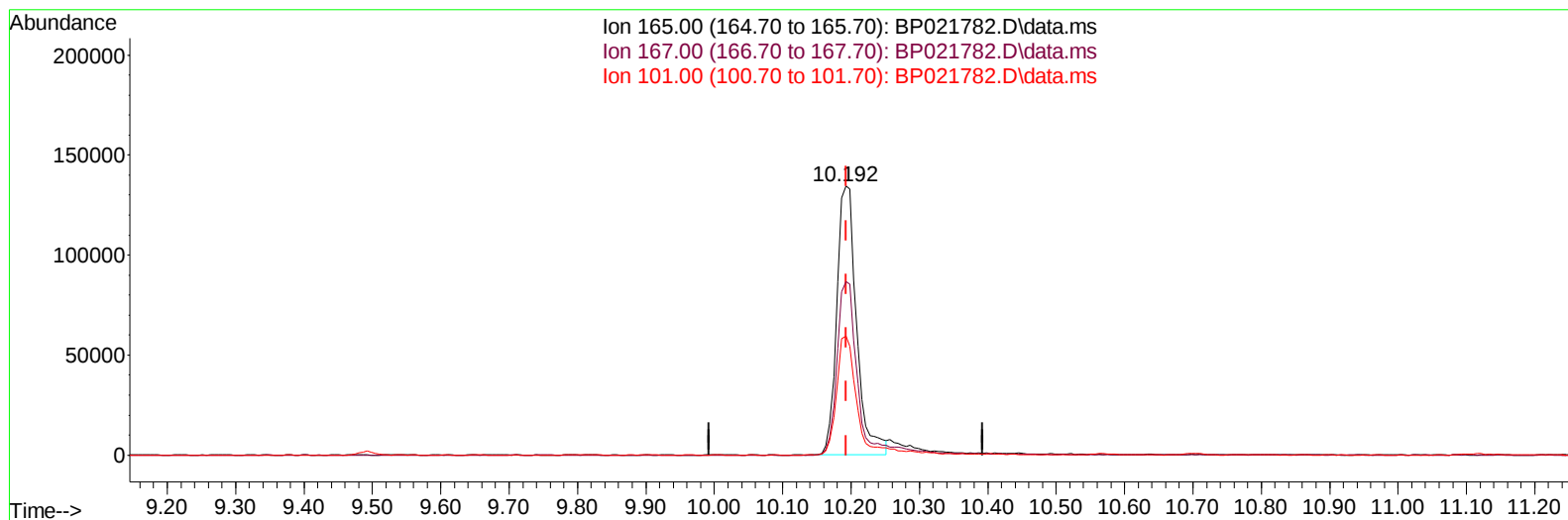
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021782.D
Acq On : 31 Aug 2024 12:36
Operator : RC/JU
Sample : P3733-15
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44A7

Manual IntegrationsAPPROVED

Quant Time: Sep 02 01:10:03 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

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



TIC: BP021782.D\data.ms

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

10.192min (-0.000) 16.73 ng/u1

response 271422

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	64.56
101.00	41.40	44.18
0.00	0.00	0.00

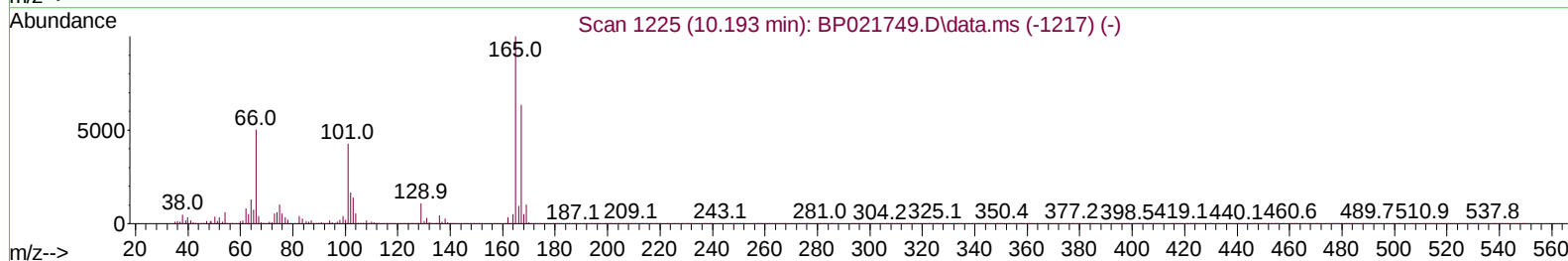
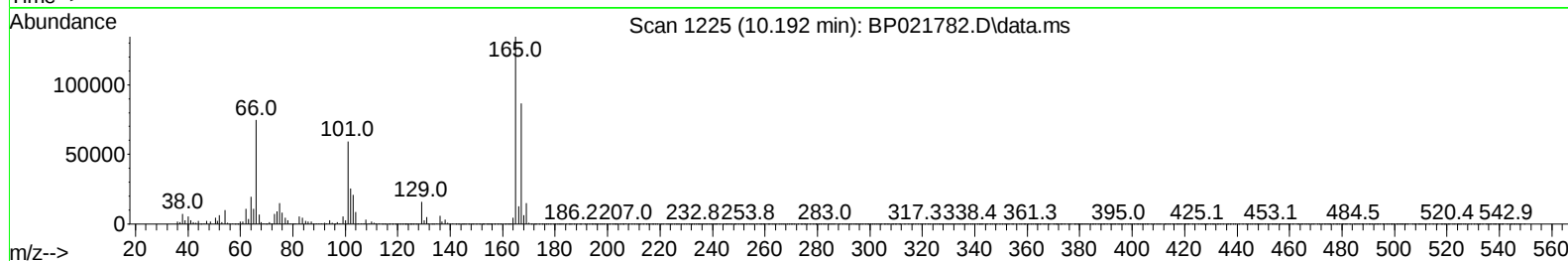
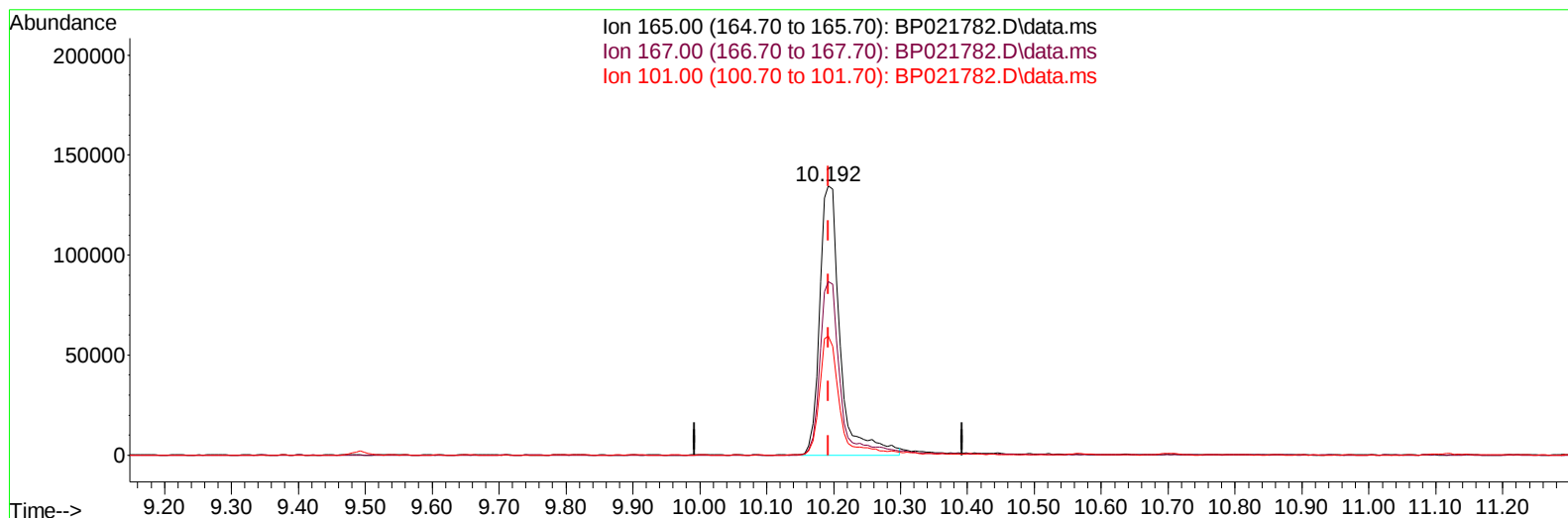
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TIC: BP021782.D\data.ms

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

10.192min (-0.000) 17.64 ng/ul m

response 286202

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	64.56
101.00	41.40	44.18
0.00	0.00	0.00

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 Sample : P3733-15
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44A7

Manual IntegrationsAPPROVED

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

Quant Time: Sep 02 01:59:32 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	252212	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	1007921	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.427	164	674091	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.221	188	1508649	20.000	ng/ul	0.00
79) Chrysene-d12	21.674	240	1514392	20.000	ng/ul	0.00
88) Perylene-d12	25.068	264	1714174	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	26505	3.489	ng/uL	0.00
4) Pyridine-d5	3.687	84	309783	14.098	ng/ul	0.00
7) Phenol-d5	6.951	99	484412	17.997	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	276278	18.861	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	325090	18.708	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113	368752	17.729	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	156348	19.280	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	152157	18.485	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.192	165	286202m	17.644	ng/ul	0.00
31) 4-Chloroaniline-d4	10.698	131	459387	19.467	ng/ul	0.00
46) Dimethylphthalate-d6	13.827	166	1085861	21.052	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	1148064	19.825	ng/ul	0.00
54) 4-Nitrophenol-d4	14.633	143	159508	16.255	ng/ul	0.00
60) Fluorene-d10	15.433	176	908687	20.955	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.545	200	112367	13.542	ng/ul	-0.01
73) Anthracene-d10	17.327	188	1520490	21.236	ng/ul	-0.01
81) Pyrene-d10	19.698	212	1851613	21.389	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.839	264	1937304	21.195	ng/ul	0.00

Target Compounds Qvalue

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

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