

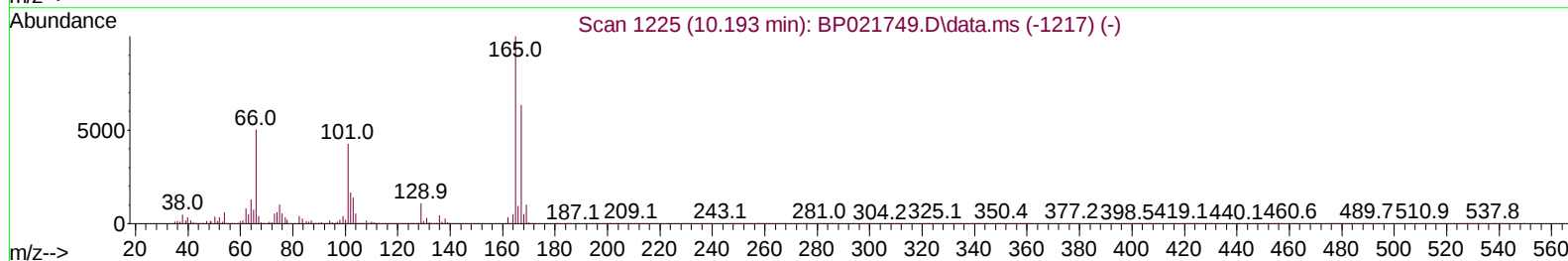
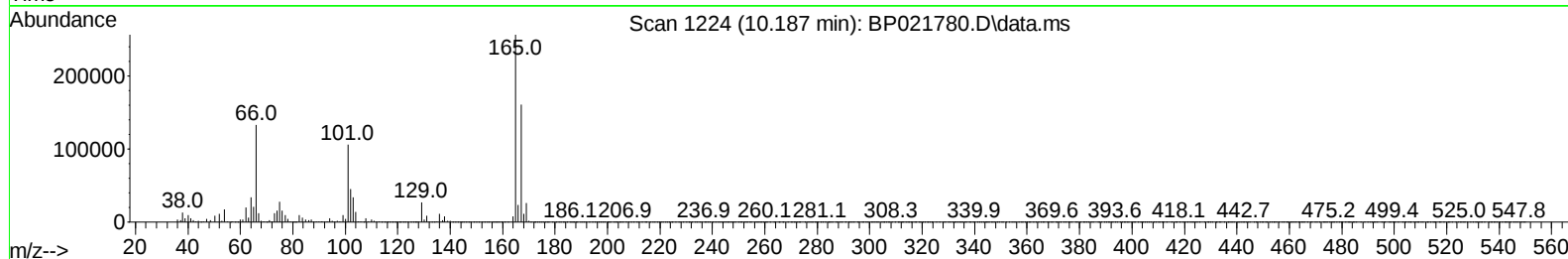
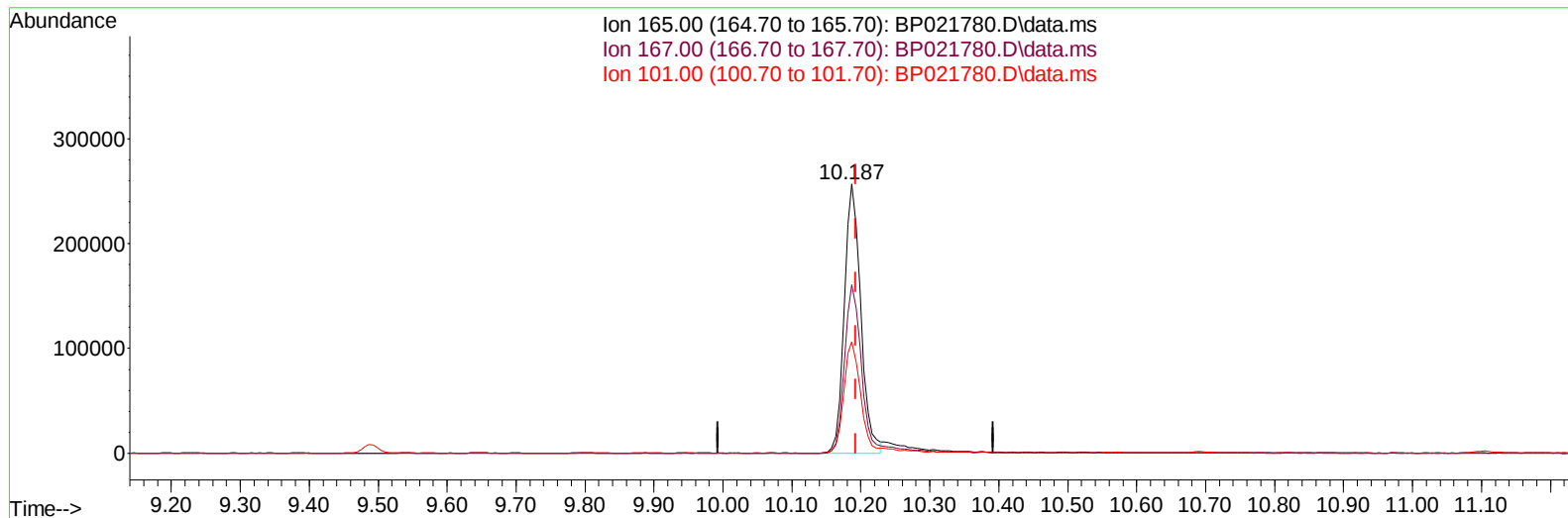
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021780.D
Acq On : 31 Aug 2024 11:09
Operator : RC/JU
Sample : P3733-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4496

Manual IntegrationsAPPROVED

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

Quant Time: Sep 02 01:09:30 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



TIC: BP021780.D\data.ms

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

10.187min (-0.006) 22.37 ng/u1

response 425123

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.63
101.00	41.40	41.48
0.00	0.00	0.00

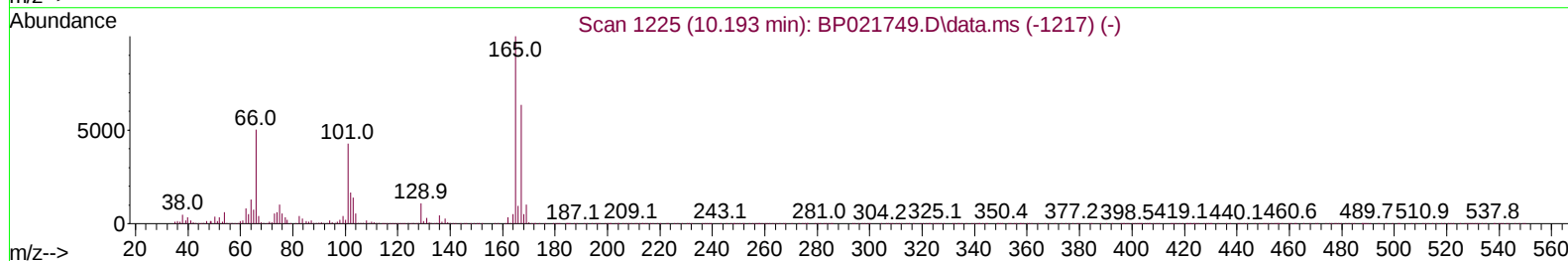
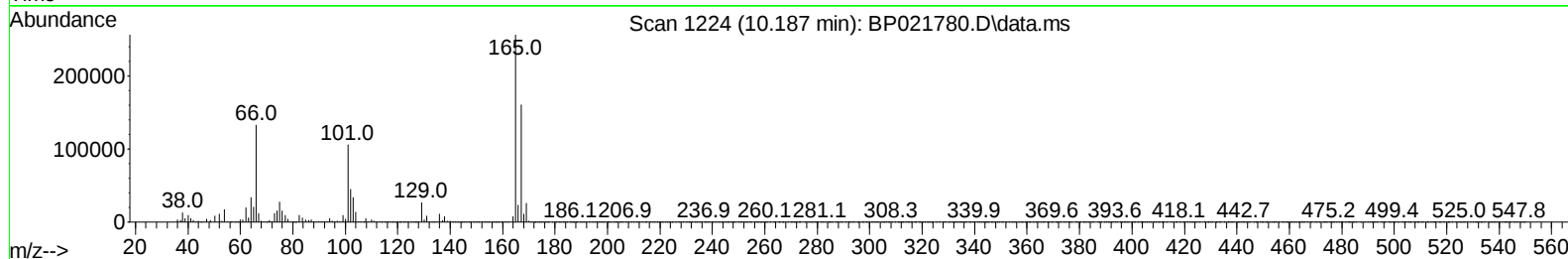
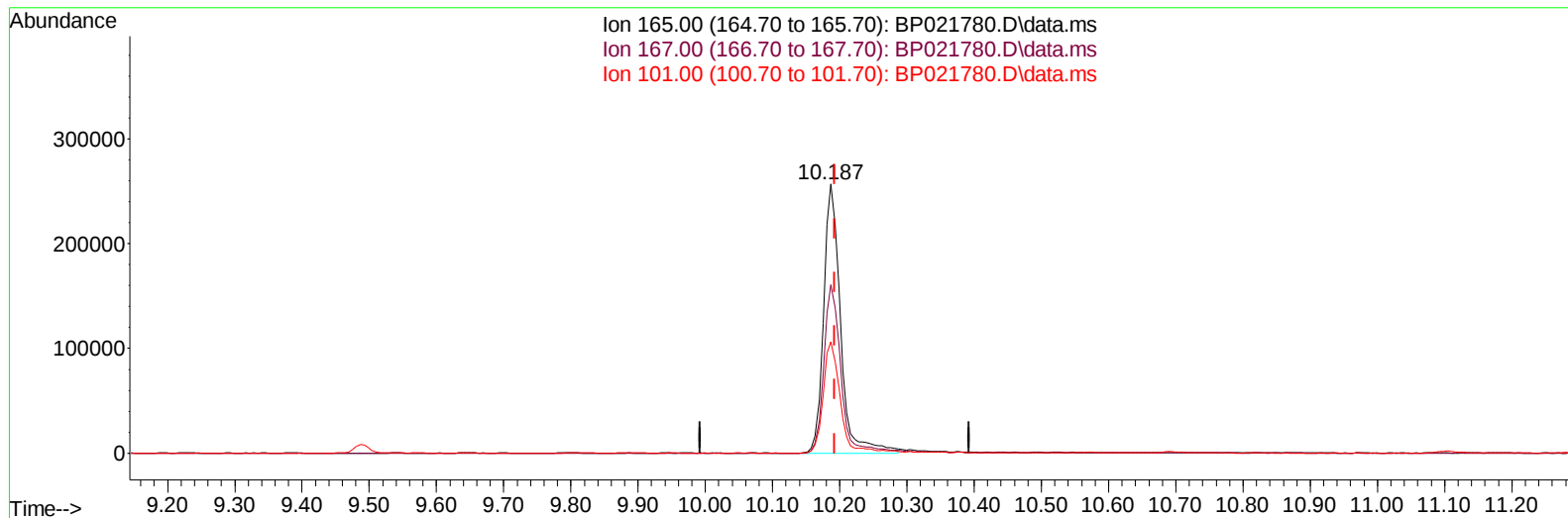
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TIC: BP021780.D\data.ms

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

10.187min (-0.006) 23.68 ng/ul m

response 449943

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.63
101.00	41.40	41.48
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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

Quant Time: Sep 02 01:50:36 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	306927	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1180907	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	689290	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.228	188	1349693	20.000	ng/ul	0.00
79) Chrysene-d12	21.686	240	1262686	20.000	ng/ul	0.02
88) Perylene-d12	25.086	264	1560162	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	36551	3.954	ng/uL	0.00
4) Pyridine-d5	3.687	84	545044	20.383	ng/ul	0.00
7) Phenol-d5	6.952	99	737512	22.516	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	402497	22.580	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	487427	23.049	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113	577493	22.816	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	237326	24.979	ng/ul	0.00
24) 2-Nitrophenol-d4	9.646	143	238361	24.716	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	449943m	23.676	ng/ul	0.00
31) 4-Chloroaniline-d4	10.699	131	671600	24.290	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1411341	26.759	ng/ul	-0.01
49) Acenaphthylene-d8	14.110	160	1610166	27.191	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.628	143	205459	20.476	ng/ul	0.00
60) Fluorene-d10	15.434	176	1207252	27.226	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.551	200	138256	18.624	ng/ul	0.00
73) Anthracene-d10	17.334	188	1824079	28.477	ng/ul	0.00
81) Pyrene-d10	19.704	212	2054364	28.462	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.857	264	2368237	28.467	ng/ul	0.02

Target Compounds Qvalue

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

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