

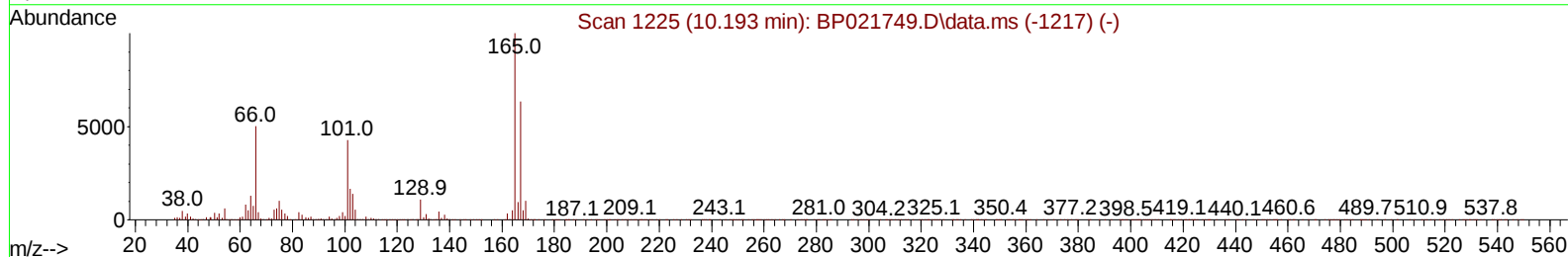
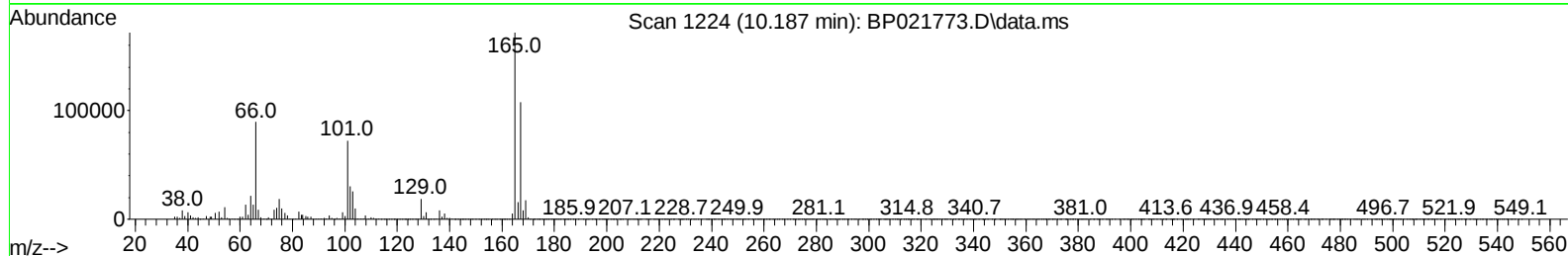
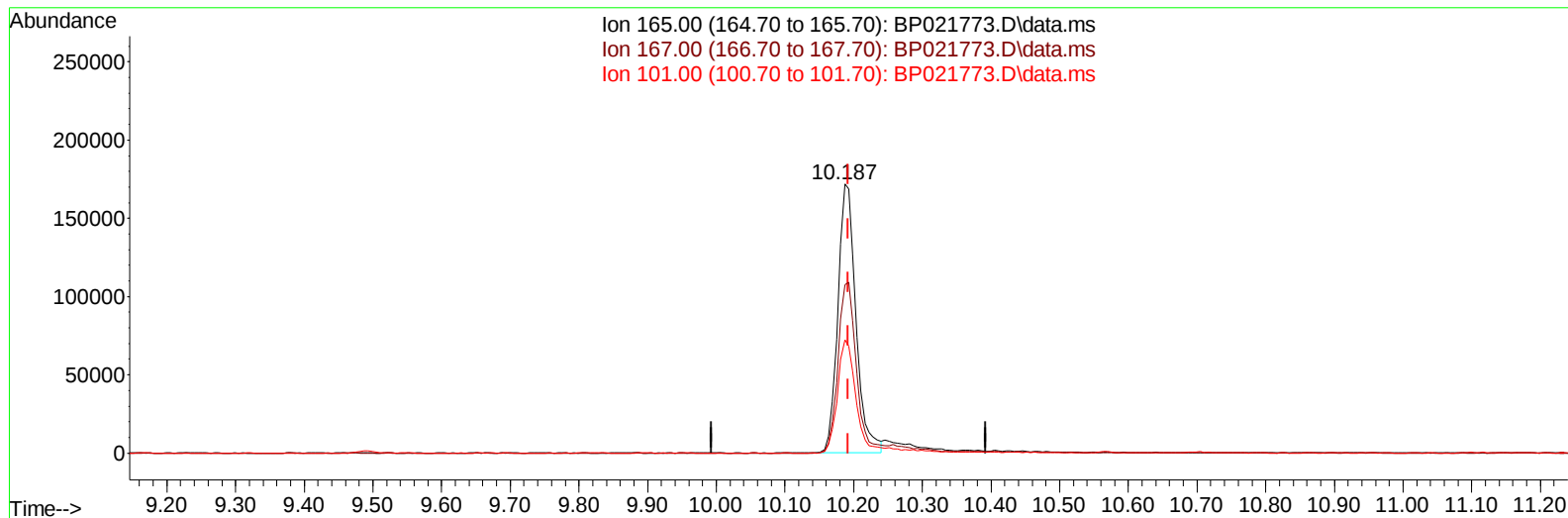
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021773.D
Acq On : 31 Aug 2024 06:06
Operator : RC/JU
Sample : P3733-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4495

Manual IntegrationsAPPROVED

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

Quant Time: Aug 31 06:35:10 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: BP021773.D\data.ms

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

10.187min (-0.006) 16.47 ng/u1

response 314573

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.64
101.00	41.40	42.09
0.00	0.00	0.00

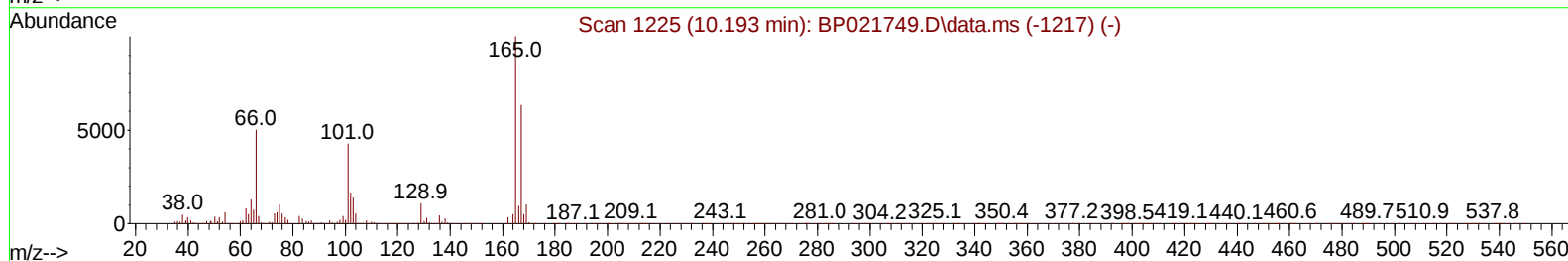
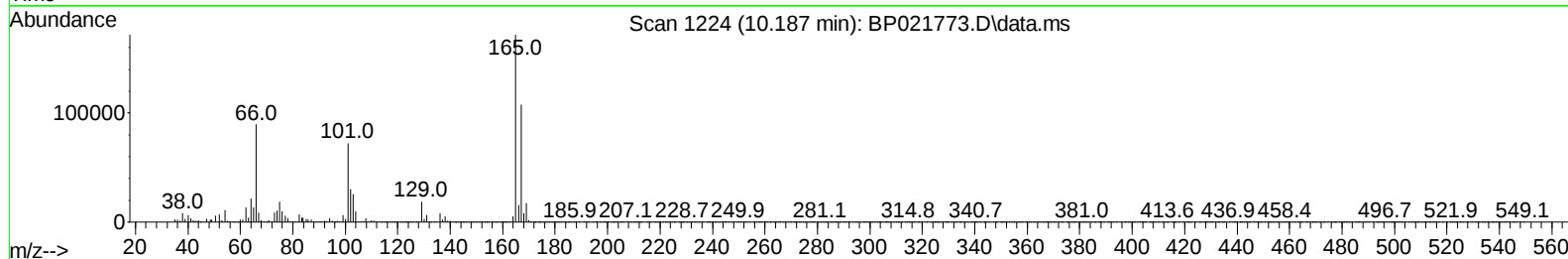
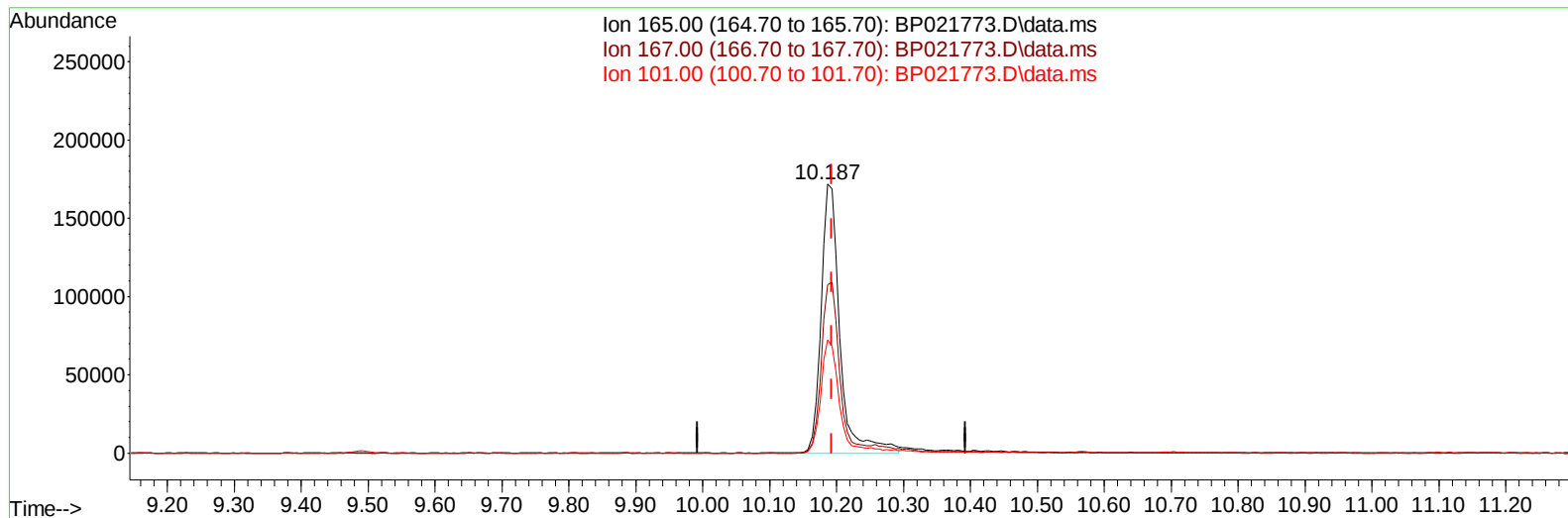
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TIC: BP021773.D\data.ms

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

10.187min (-0.006) 17.52 ng/ul m

response 334655

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.64
101.00	41.40	42.09
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/02/2024
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Quant Time: Aug 31 06:36:11 2024
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 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	320982	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1186620	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	696193	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1466737	20.000	ng/ul	0.00
79) Chrysene-d12	21.675	240	1581988	20.000	ng/ul	0.00
88) Perylene-d12	25.069	264	1892035	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	35769	3.700	ng/uL	0.00
4) Pyridine-d5	3.687	84	482451	17.252	ng/ul	0.00
7) Phenol-d5	6.952	99	619159	18.075	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	346192	18.571	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	418897	18.941	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	445754	16.840	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	186510	19.536	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	180577	18.634	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	334655m	17.524	ng/ul	0.00
31) 4-Chloroaniline-d4	10.699	131	496098	17.856	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	983671	18.465	ng/ul	-0.01
49) Acenaphthylene-d8	14.110	160	1114724	18.638	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.628	143	143481	14.157	ng/ul	0.00
60) Fluorene-d10	15.434	176	862709	19.263	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.551	200	93309	11.567	ng/ul	0.00
73) Anthracene-d10	17.334	188	1361495	19.559	ng/ul	0.00
81) Pyrene-d10	19.698	212	1708971	18.898	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.839	264	1953841	19.366	ng/ul	0.00

Target Compounds Qvalue

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

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