

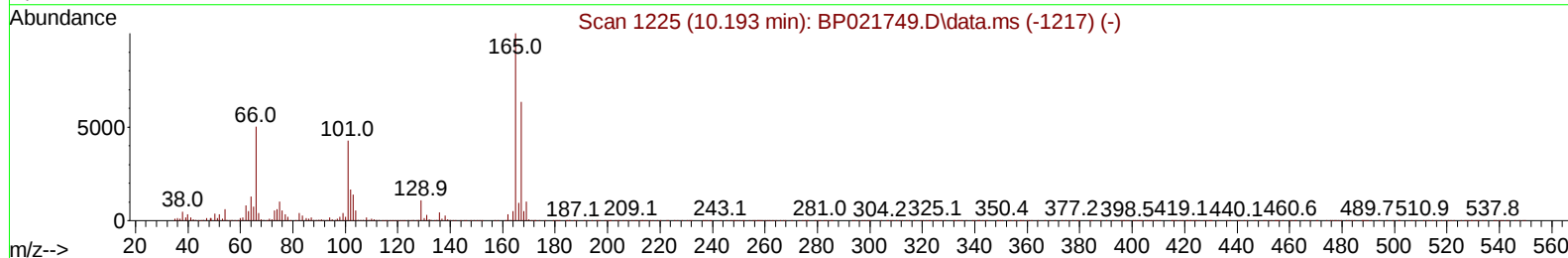
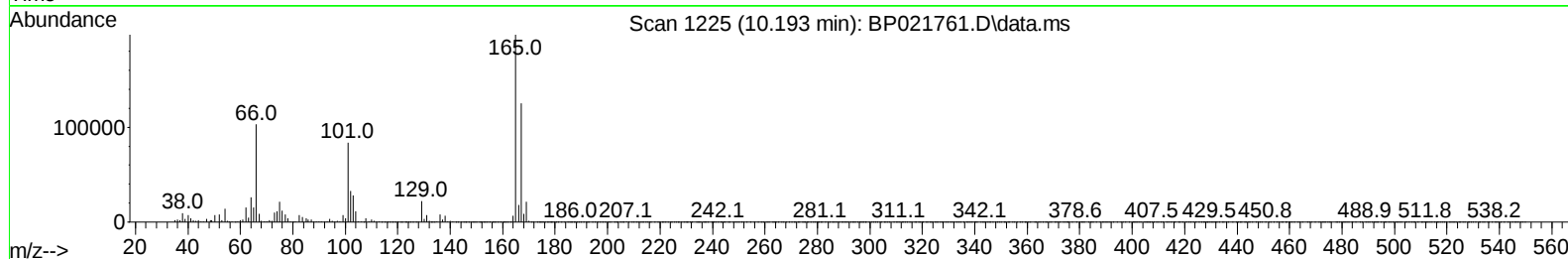
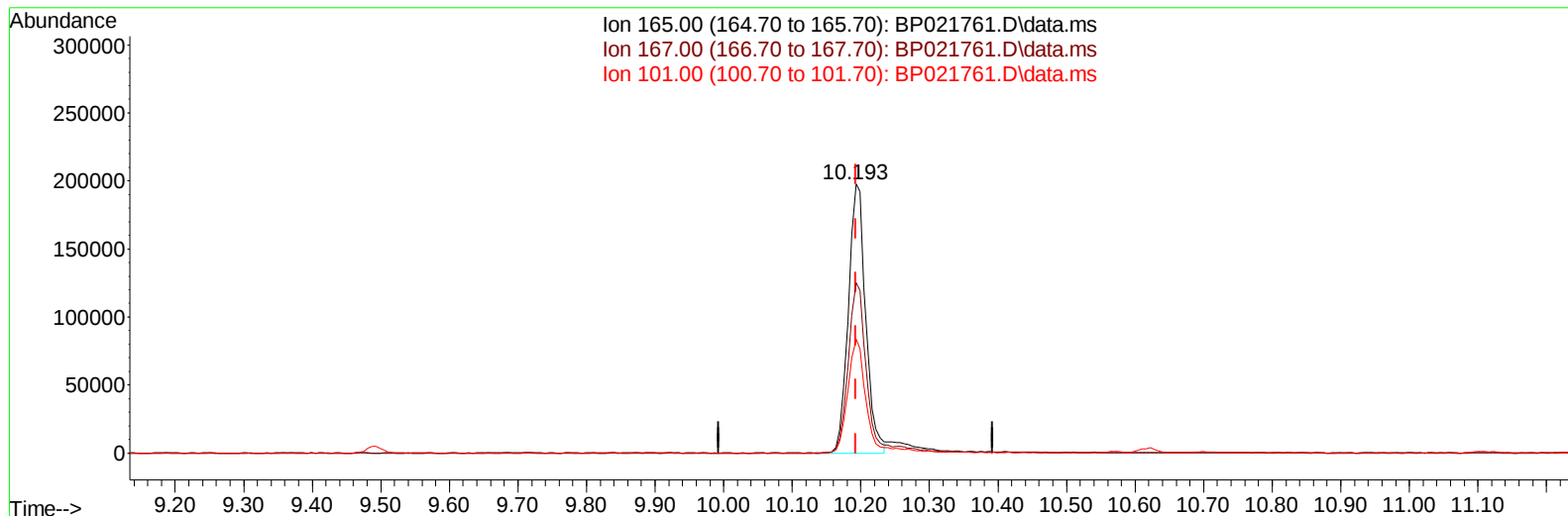
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021761.D
Acq On : 30 Aug 2024 20:36
Operator : RC/JU
Sample : P3438-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYL3

Manual IntegrationsAPPROVED

Quant Time: Aug 30 23:09:49 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 09/02/2024
Supervised By :mohammad ahmed 09/03/2024



TIC: BP021761.D\data.ms

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

10.193min (-0.000) 13.42 ng/u1

response 346207

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	63.33
101.00	41.40	42.40
0.00	0.00	0.00

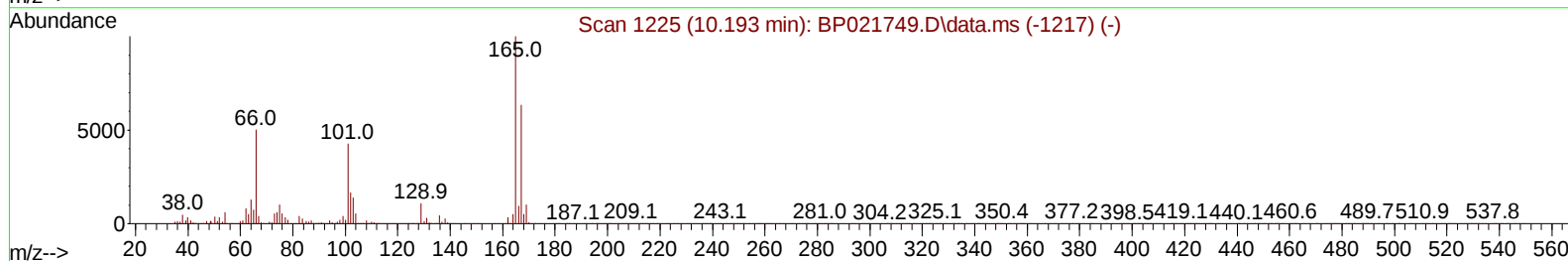
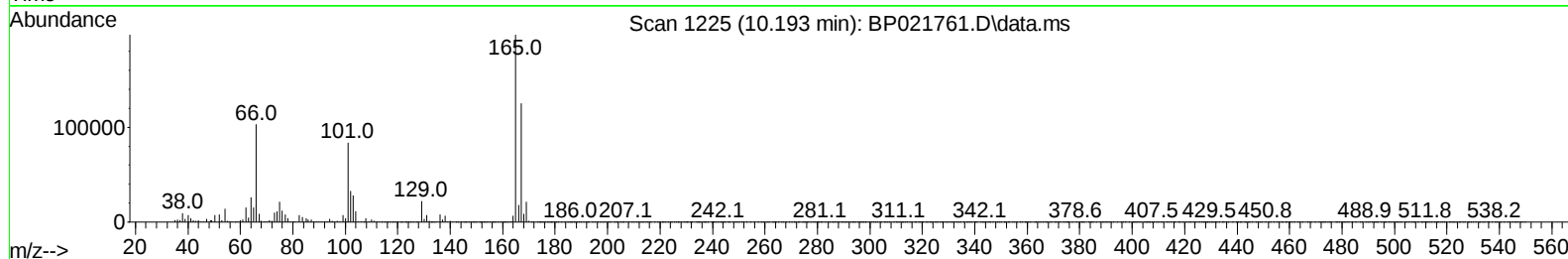
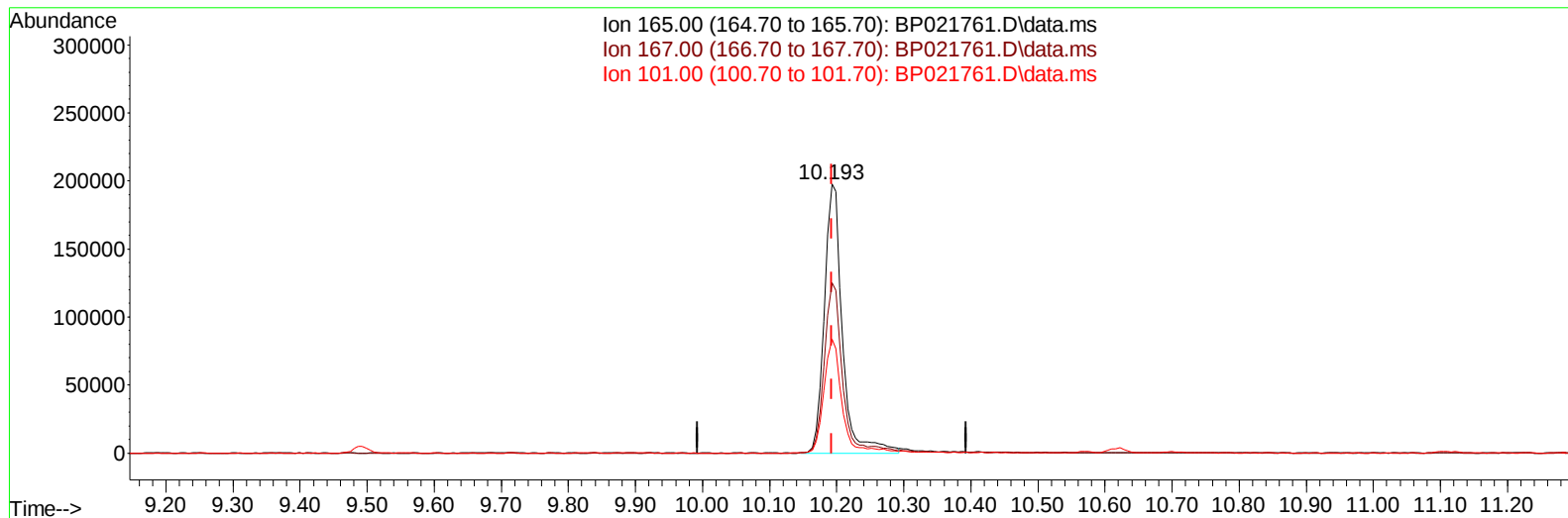
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TIC: BP021761.D\data.ms

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

10.193min (-0.000) 14.26 ng/ul m

response 367853

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	63.33
101.00	41.40	42.40
0.00	0.00	0.00

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 BNA_P
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 DCYL3

Manual IntegrationsAPPROVED

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

Quant Time: Aug 30 23:26:39 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	403171	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1602803	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	1016955	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	2075162	20.000	ng/ul	0.00
79) Chrysene-d12	21.686	240	2047146	20.000	ng/ul	0.02
88) Perylene-d12	25.092	264	2420815	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	32040	2.638	ng/uL	0.00
4) Pyridine-d5	3.687	84	417145	11.876	ng/ul	0.00
7) Phenol-d5	6.958	99	600535	13.957	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	335862	14.344	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	408393	14.702	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	466622	14.035	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	196081	15.206	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	193985	14.820	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	367853m	14.261	ng/ul	0.00
31) 4-Chloroaniline-d4	10.699	131	522790	13.931	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1193020	15.332	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	1367812	15.656	ng/ul	0.00
54) 4-Nitrophenol-d4	14.640	143	169775	11.468	ng/ul	0.01
60) Fluorene-d10	15.440	176	1049517	16.043	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	114633	10.044	ng/ul	0.00
73) Anthracene-d10	17.328	188	1612548	16.374	ng/ul	-0.01
81) Pyrene-d10	19.704	212	1864436	15.932	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.857	264	2123772	16.452	ng/ul	0.02
Target Compounds						
					Qvalue	
30) Naphthalene	10.616	128	191252	2.117	ng/ul	99
52) Acenaphthene	14.493	153	70836	1.072	ng/ul	99
72) Phenanthrene	17.269	178	211691	1.858	ng/ul	99

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

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