

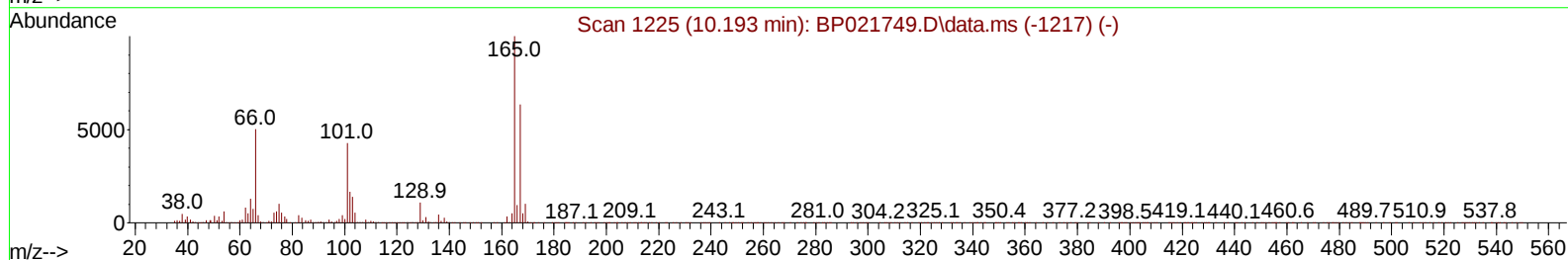
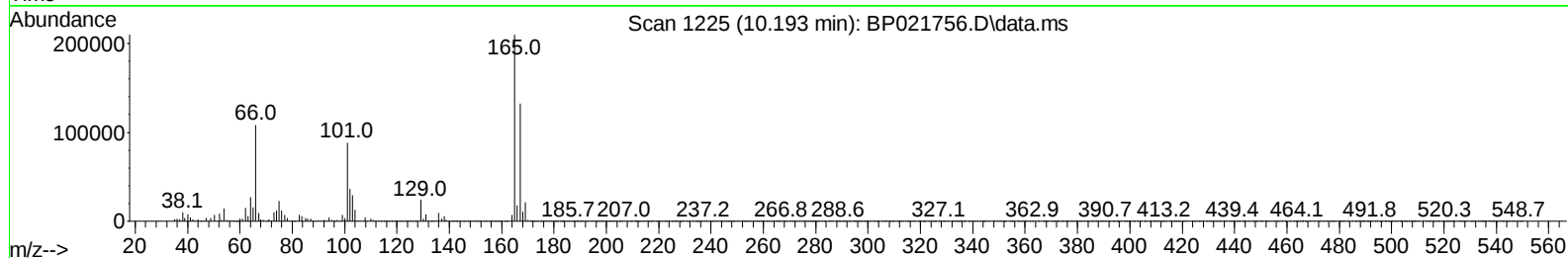
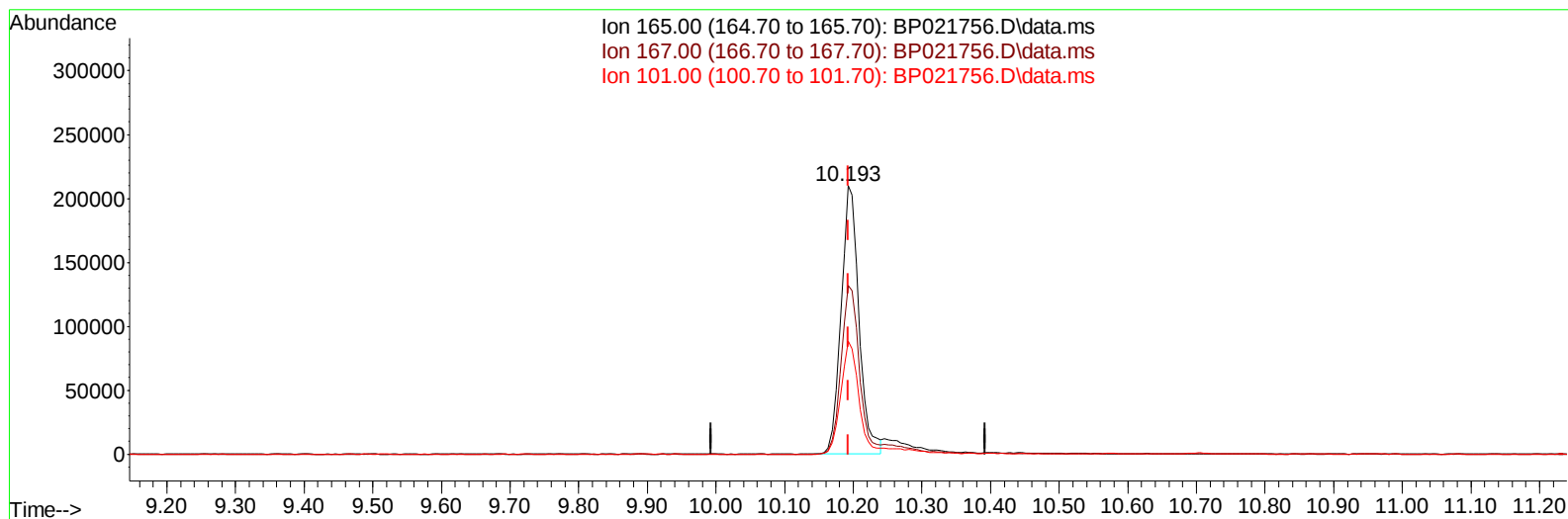
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
 Data File : BP021756.D
 Acq On : 30 Aug 2024 17:00
 Operator : RC/JU
 Sample : PB163016BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK016

Manual IntegrationsAPPROVED

Quant Time: Aug 30 17:45:21 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: BP021756.D\data.ms

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

10.193min (-0.000) 25.94 ng/u1

response 380328

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.91
101.00	41.40	42.13
0.00	0.00	0.00

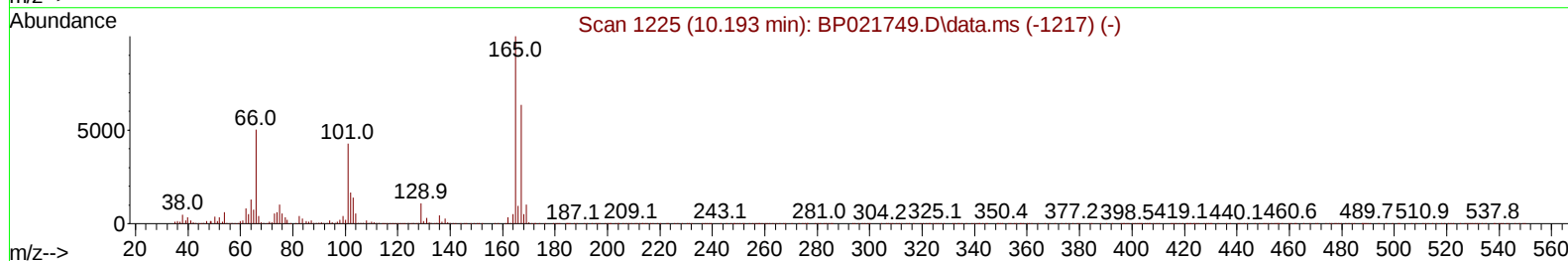
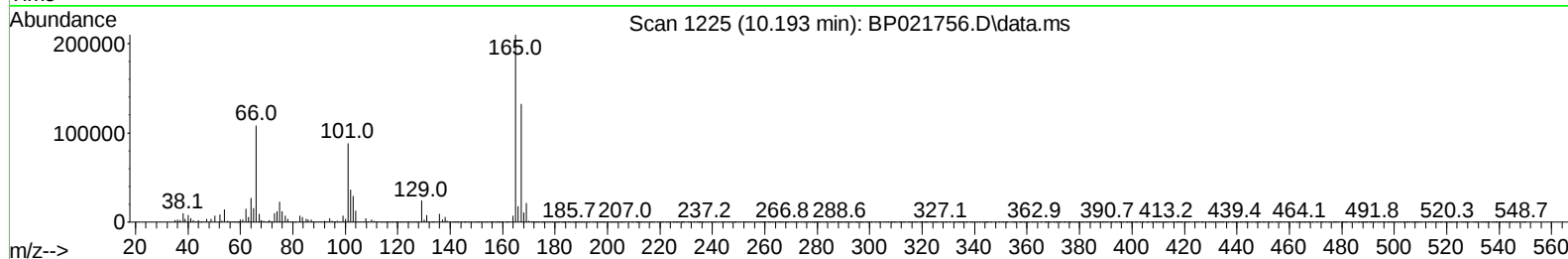
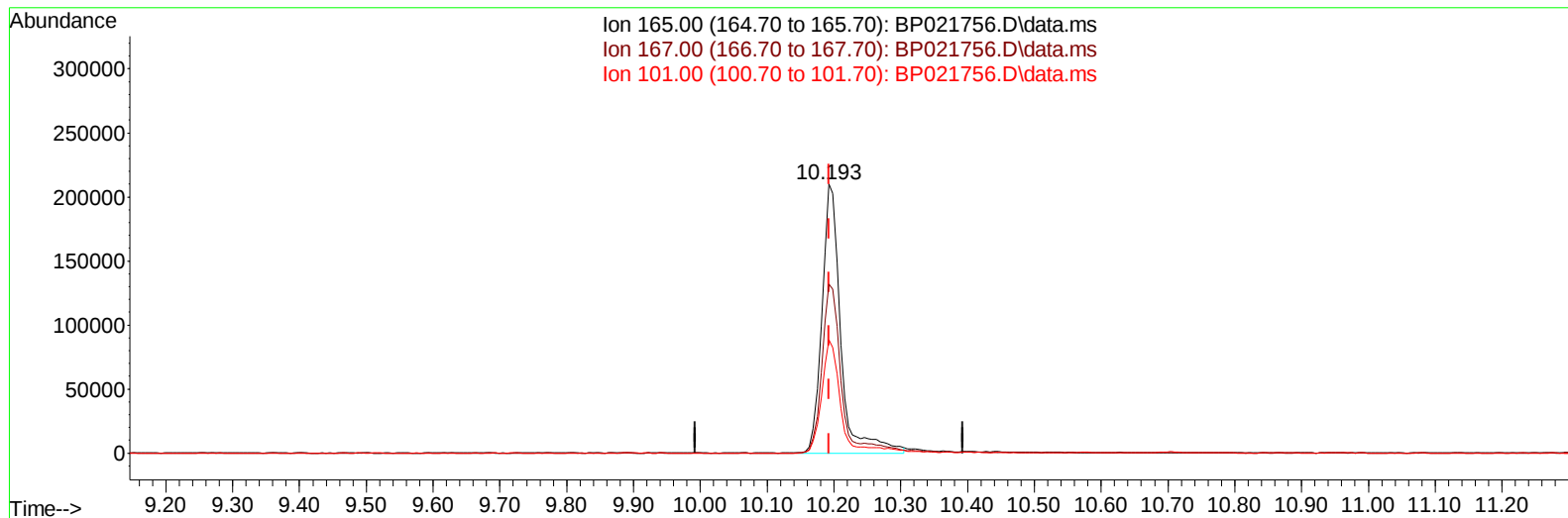
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TIC: BP021756.D\data.ms

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

10.193min (-0.000) 28.11 ng/ul m

response 412167

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.91
101.00	41.40	42.13
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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

Quant Time: Aug 31 00:01:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	230005	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	911085	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.434	164	597856	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	1363134	20.000	ng/ul	0.01
79) Chrysene-d12	21.680	240	1405634	20.000	ng/ul	0.01
88) Perylene-d12	25.092	264	1544894	20.000	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	44910	6.482	ng/uL	0.00
4) Pyridine-d5	3.687	84	576071	28.748	ng/ul	0.00
7) Phenol-d5	6.957	99	709581	28.908	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	385981	28.895	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	468477	29.562	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	520319	27.432	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	227191	30.994	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	237436	31.912	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	412167m	28.111	ng/ul	0.00
31) 4-Chloroaniline-d4	10.704	131	588534	27.590	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1431588	31.294	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	1540052	29.985	ng/ul	0.00
54) 4-Nitrophenol-d4	14.645	143	245289	28.183	ng/ul	0.02
60) Fluorene-d10	15.439	176	1205738	31.351	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	183377	24.459	ng/ul	0.00
73) Anthracene-d10	17.333	188	1989200	30.748	ng/ul	0.00
81) Pyrene-d10	19.704	212	2514978	31.300	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.857	264	2549956	30.954	ng/ul	0.02

Target Compounds Qvalue

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

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