

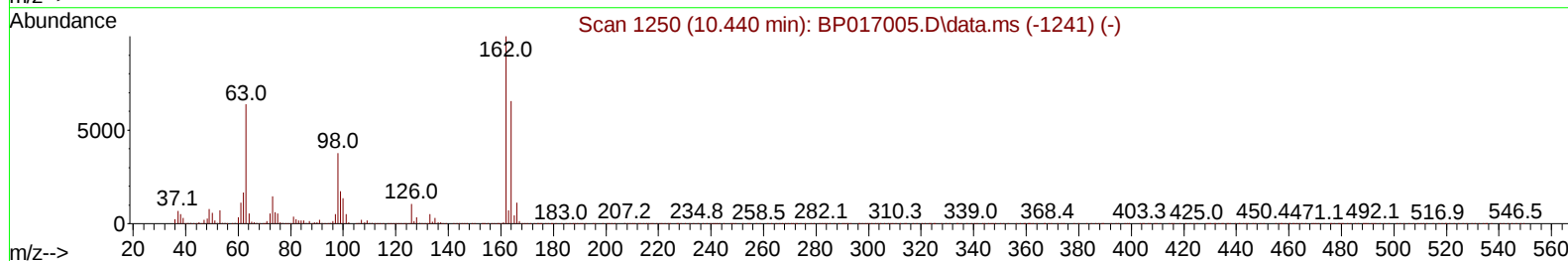
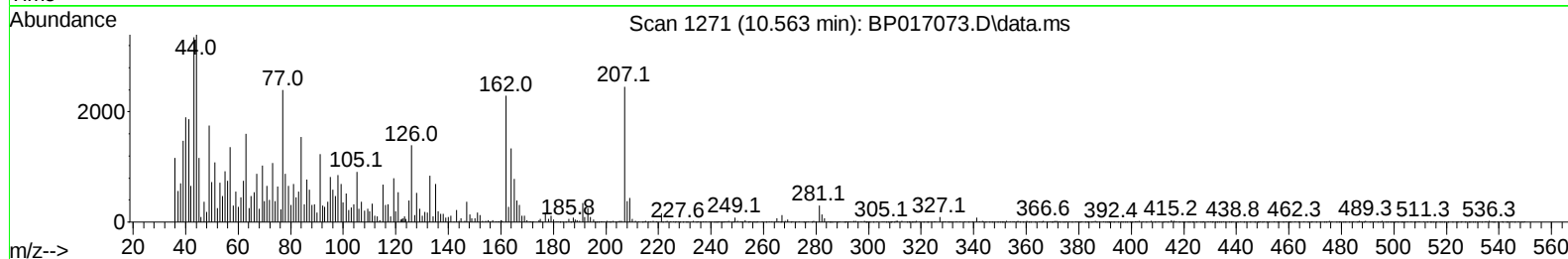
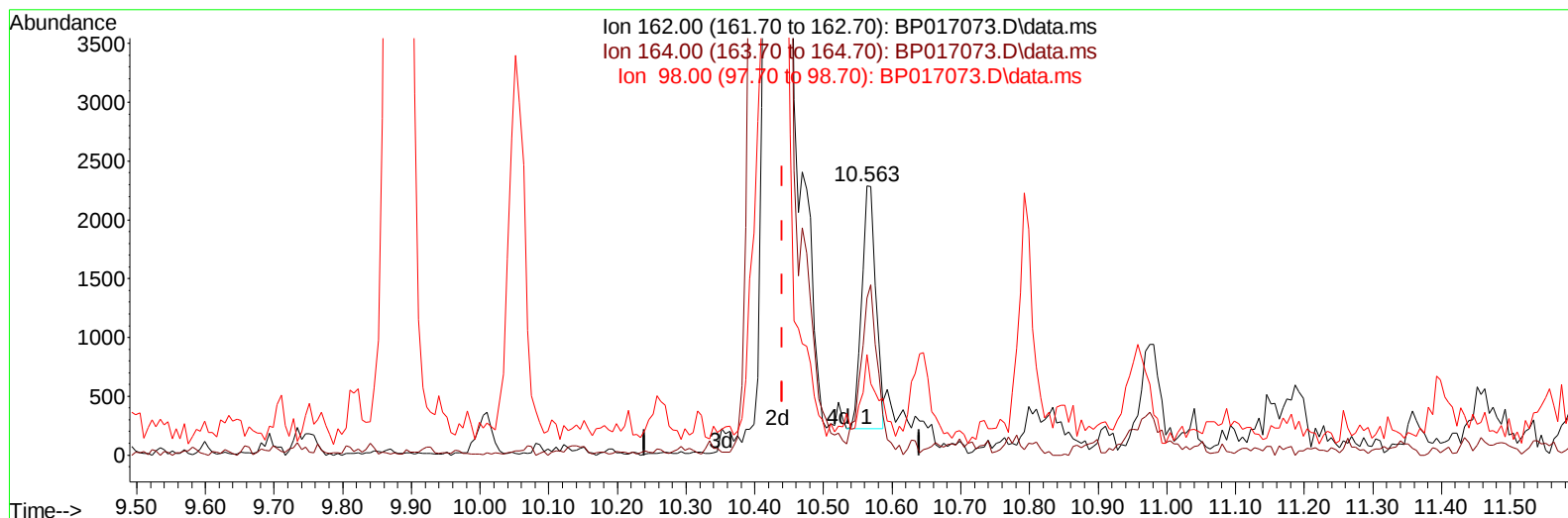
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017073.D
 Acq On : 10 Sep 2023 03:49
 Operator : MA/JU
 Sample : 04227-10
 Misc :
 ALS Vial : 85 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKE0

Manual IntegrationsAPPROVED

Quant Time: Sep 10 04:56:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/11/2023



TIC: BP017073.D\data.ms

(29) 2,4-Dichlorophenol (C)

10.563min (+ 0.123) 0.10 ng/u1

response 2990

Ion	Exp%	Act%
162.00	100.00	100.00
164.00	64.70	58.42
98.00	35.30	37.30
0.00	0.00	0.00

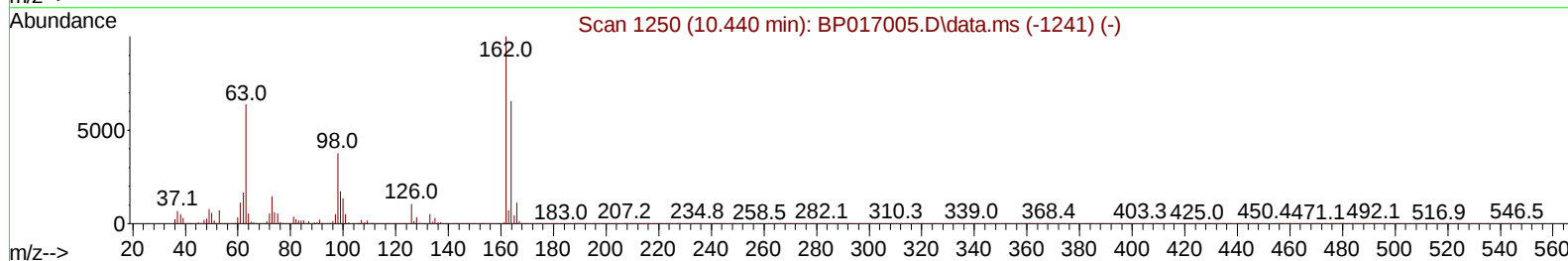
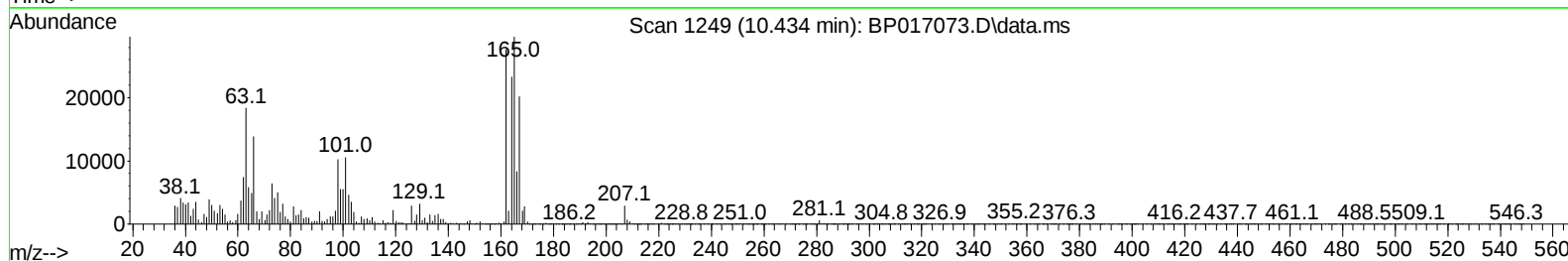
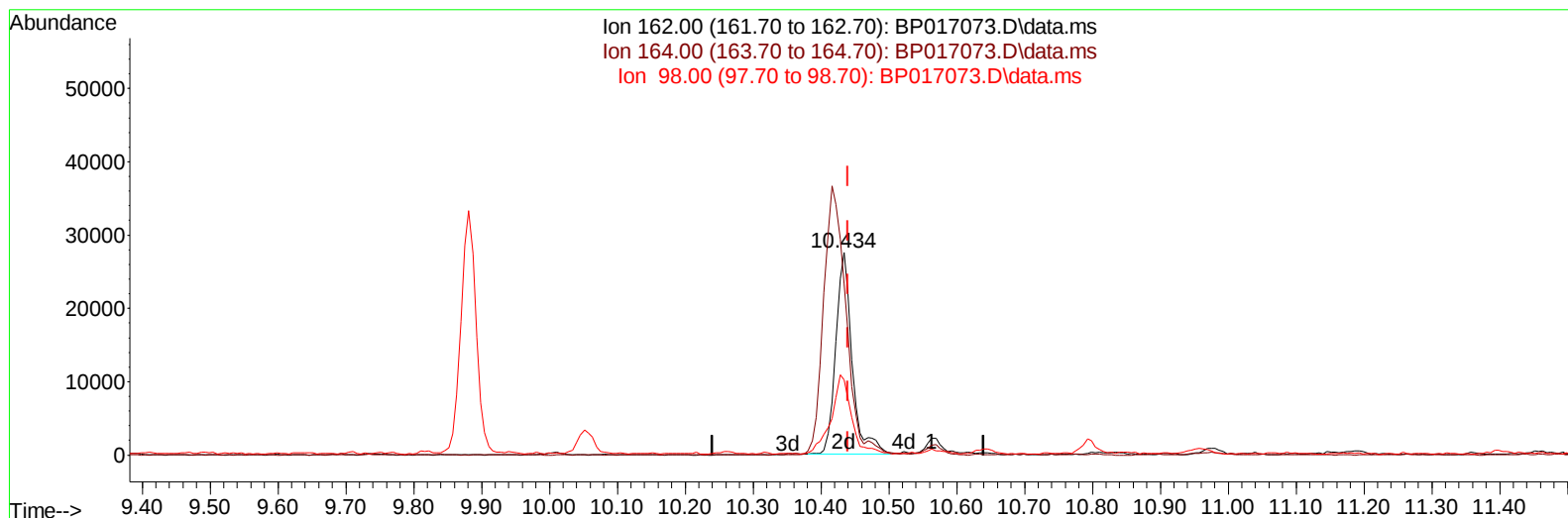
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TIC: BP017073.D\data.ms

(29) 2,4-Dichlorophenol (C)

10.434min (-0.006) 1.51 ng/ul m

response 46575

Ion	Exp%	Act%
162.00	100.00	100.00
164.00	64.70	84.58#
98.00	35.30	37.19
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/11/2023

Quant Time: Sep 10 04:59:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.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	7.969	152	443286	20.000	ng/ul	0.00
20) Naphthalene-d8	10.793	136	1945101	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.628	164	1143610	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	2304842	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1517222	20.000	ng/ul	0.00
88) Perylene-d12	24.021	264	1417310	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	62360	5.321	ng/uL	0.00
4) Pyridine-d5	3.764	84	403514	12.050	ng/ul	0.00
7) Phenol-d5	7.116	99	358049	8.841	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	1114151	43.982	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	1003521	34.238	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113	684063	22.172	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	641703	42.524	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	680196	41.582	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	1152841	38.174	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	1358024	31.168	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	3960147	46.520	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	4480194	45.604	ng/ul	0.00
54) 4-Nitrophenol-d4	14.845	143	113908	6.850	ng/ul	0.00
60) Fluorene-d10	15.622	176	3352878	46.002	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	507682	36.602	ng/ul	0.00
73) Anthracene-d10	17.498	188	5032523	47.545	ng/ul	0.00
81) Pyrene-d10	19.733	212	5046989	54.613	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.857	264	3523974	49.010	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.358	88	34761	2.747	ng/uL	97
29) 2,4-Dichlorophenol	10.434	162	46575m	1.513	ng/ul	
42) 2,4,5-Trichlorophenol	13.128	196	46308	1.856	ng/ul	94

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

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