

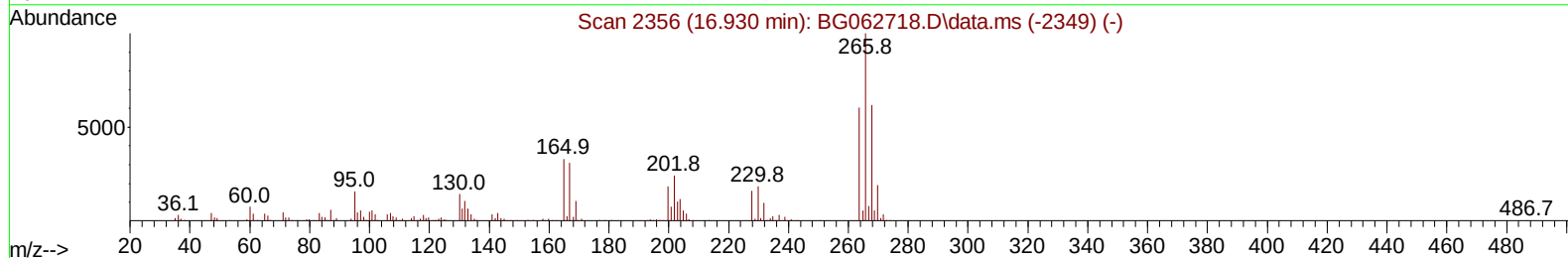
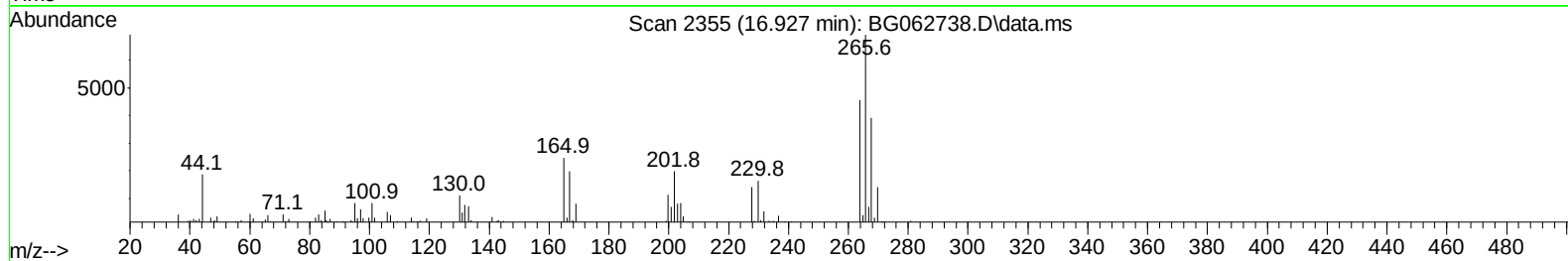
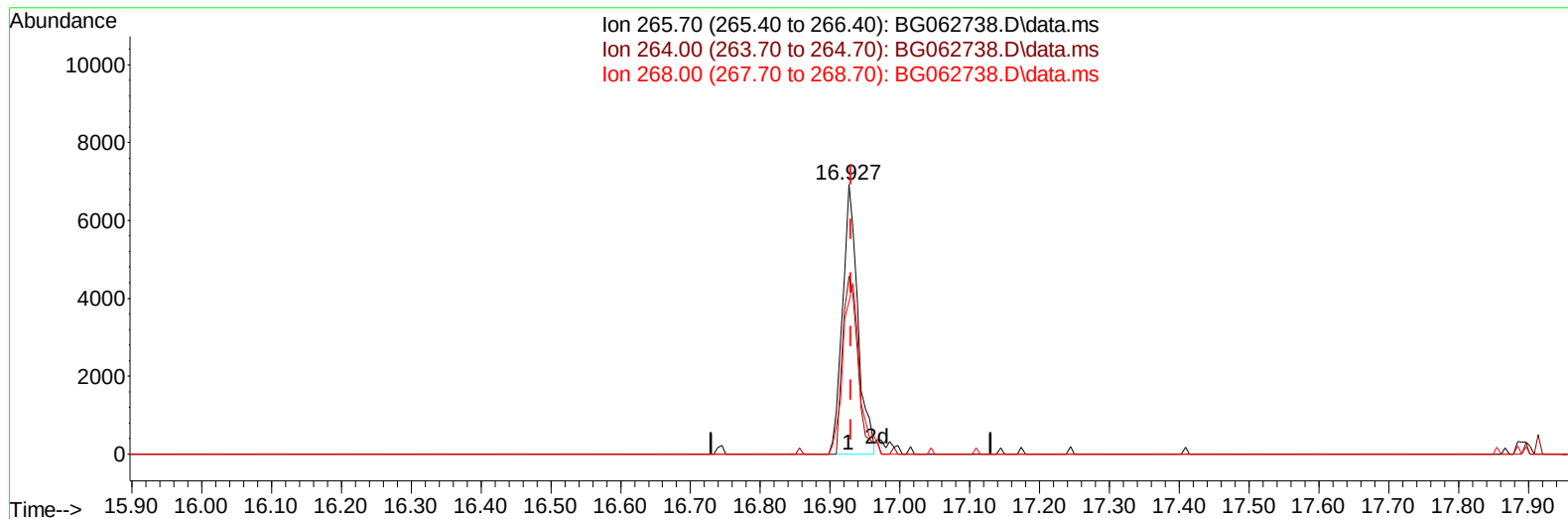
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091024\
Data File : BG062738.D
Acq On : 11 Sep 2024 16:27
Operator : JU/RC
Sample : P3877-17
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ6

Manual IntegrationsAPPROVED

Quant Time: Sep 11 23:04:55 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/12/2024
Supervised By :mohammad ahmed 09/13/2024



TIC: BG062738.D\data.ms

(71) Pentachlorophenol (C)

16.927min (-0.003) 1.72 ng/u1

response 10470

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	60.70	65.84
268.00	61.90	56.42
0.00	0.00	0.00

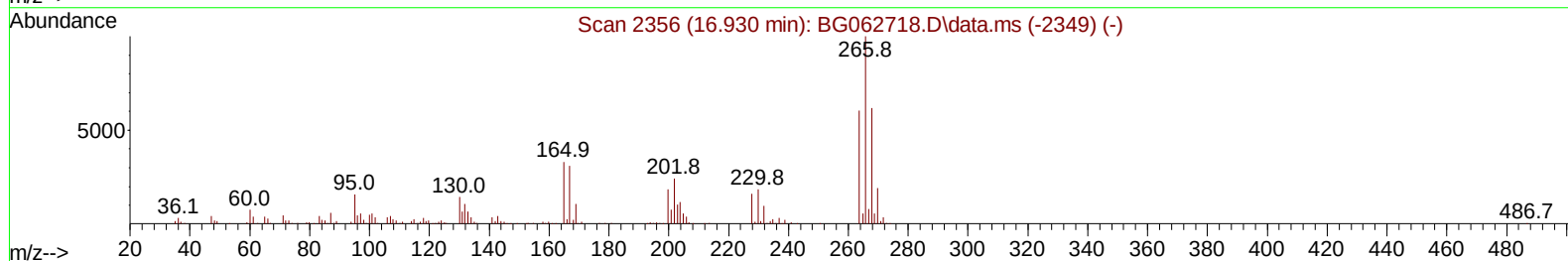
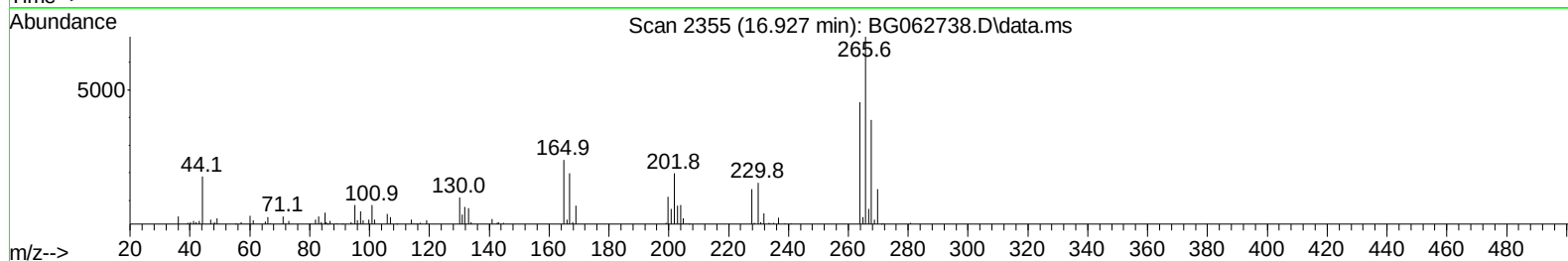
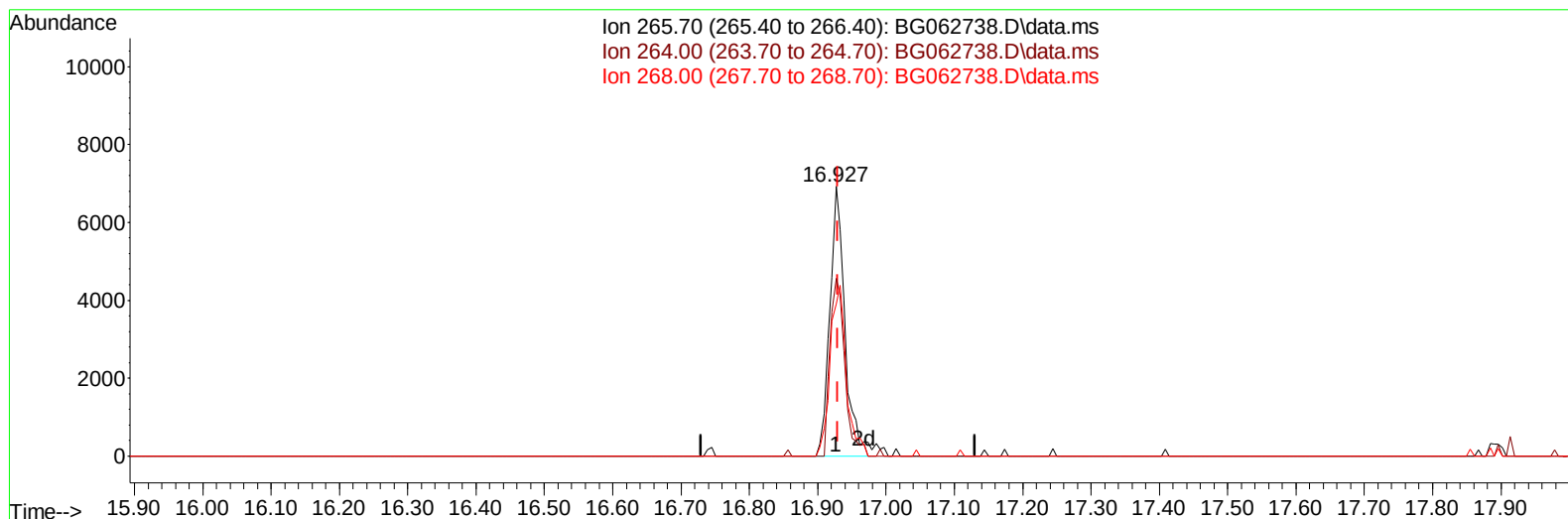
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TIC: BG062738.D\data.ms

(71) Pentachlorophenol (C)

16.927min (-0.003) 1.81 ng/ul m

response 11028

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	60.70	65.84
268.00	61.90	56.42
0.00	0.00	0.00

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Quant Time: Sep 11 23:40:01 2024
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 QLast Update : Tue Sep 10 01:28:17 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.902	152	82161	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	339858	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.530	164	287764	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.273	188	775761	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	805733	20.000	ng/ul	0.00
88) Perylene-d12	24.588	264	875249	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	4250	2.338	ng/uL	0.00
4) Pyridine-d5	3.766	84	68249	11.796	ng/ul	0.00
7) Phenol-d5	7.068	99	95235	13.219	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.232	67	54487	12.614	ng/ul	0.00
11) 2-Chlorophenol-d4	7.432	132	71888	13.699	ng/ul	0.00
15) 4-Methylphenol-d8	8.601	113	82498	13.741	ng/ul	0.00
21) Nitrobenzene-d5	9.059	128	34524	14.213	ng/ul	0.00
24) 2-Nitrophenol-d4	9.776	143	42952	16.169	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.317	165	87006	15.039	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	111080	13.968	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	321413	14.504	ng/ul	0.00
49) Acenaphthylene-d8	14.224	160	358631	14.572	ng/ul	0.00
54) 4-Nitrophenol-d4	14.723	143	43060	11.439	ng/ul	0.00
60) Fluorene-d10	15.523	176	310975	15.573	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	44160	9.896	ng/ul	0.00
73) Anthracene-d10	17.373	188	573600	15.667	ng/ul	0.00
81) Pyrene-d10	19.665	212	741655	16.357	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.377	264	741903	16.053	ng/ul	0.00
Target Compounds						
71) Pentachlorophenol	16.927	266	11028m	1.809	ng/ul	Qvalue

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

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