

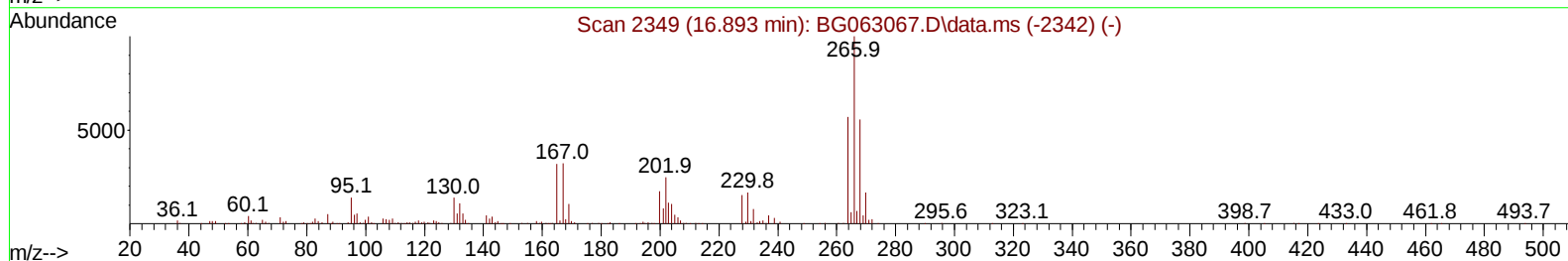
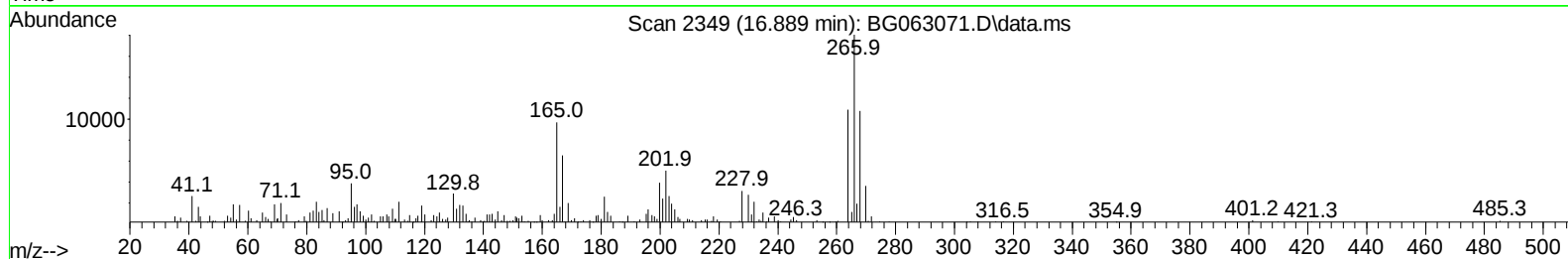
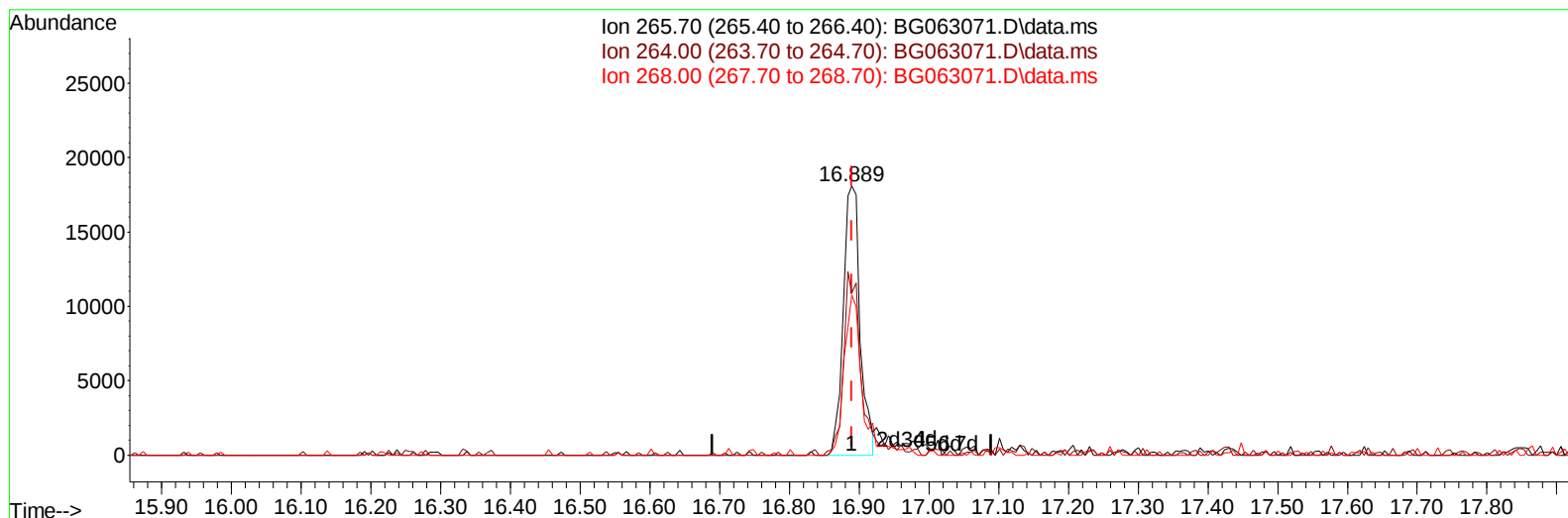
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCC7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/30/2024
Supervised By :mohammad ahmed 10/01/2024

Quant Time: Sep 27 22:55:13 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 25 13:16:45 2024
Response via : Initial Calibration



TIC: BG063071.D\data.ms

(71) Pentachlorophenol (C)

16.889min (-0.000) 4.71 ng/u1

response 30176

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	72.10	60.35
268.00	71.60	59.60
0.00	0.00	0.00

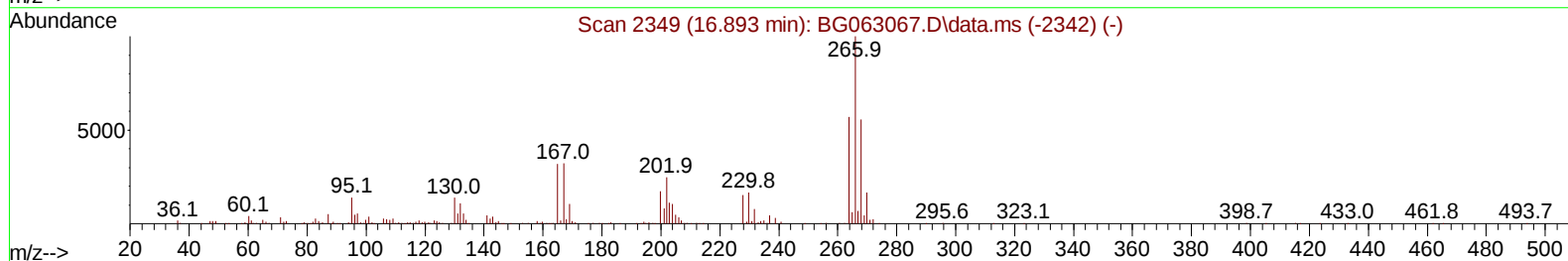
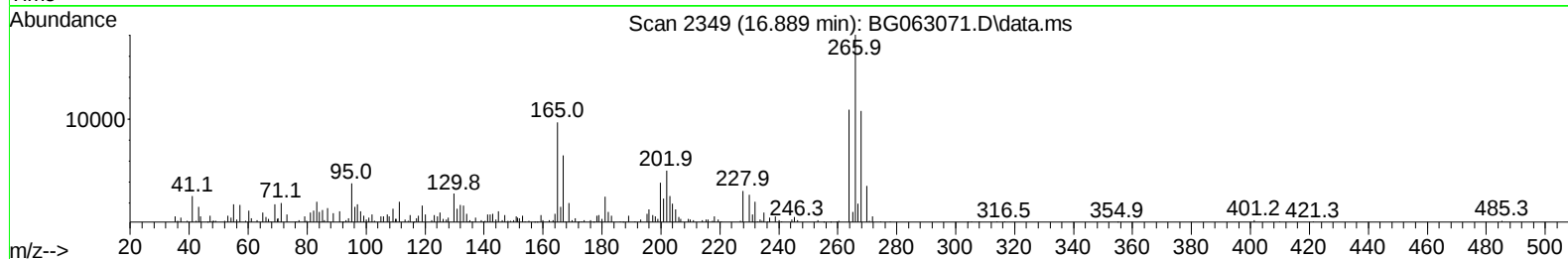
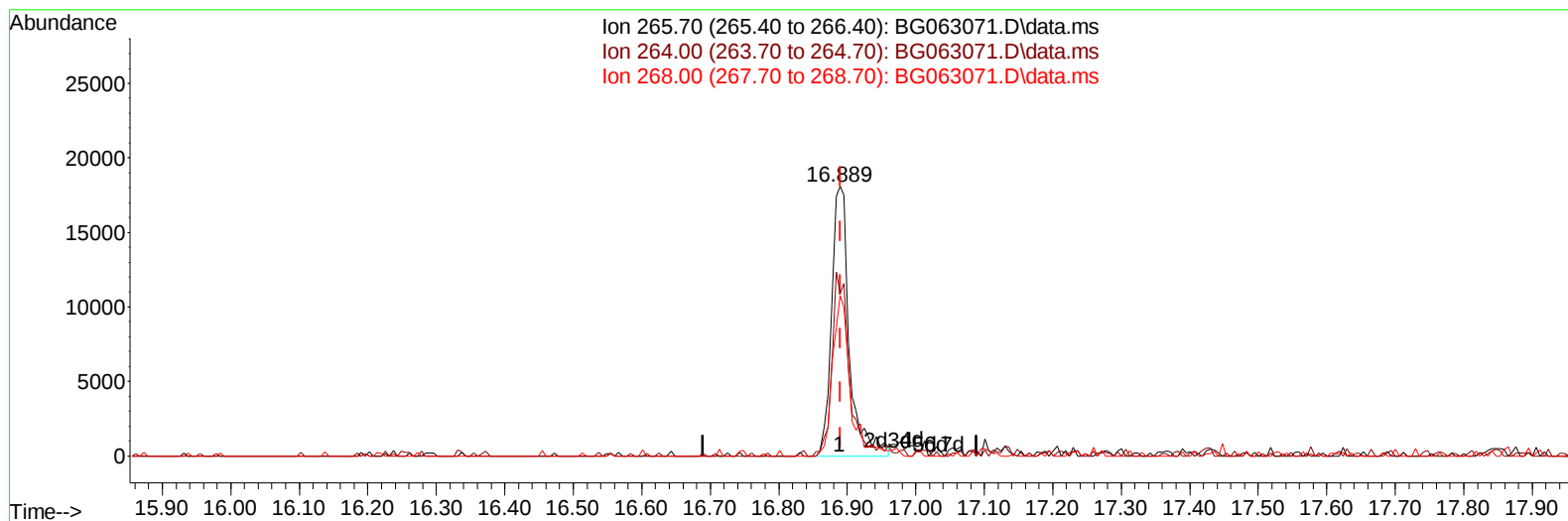
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TIC: BG063071.D\data.ms

(71) Pentachlorophenol (C)

16.889min (-0.000) 5.11 ng/ul m

response 32734

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	72.10	60.35
268.00	71.60	59.60
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/30/2024
 Supervised By :mohammad ahmed 10/01/2024

Quant Time: Sep 27 23:22:53 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 25 13:16:45 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.847	152	74190	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.638	136	307391	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.480	164	274311	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.236	188	753365	20.000	ng/ul	0.00
79) Chrysene-d12	21.484	240	802176	20.000	ng/ul	# 0.00
88) Perylene-d12	24.522	264	944175	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	3945	3.065	ng/uL	0.00
4) Pyridine-d5	3.722	84	43965	13.971	ng/ul	0.00
7) Phenol-d5	7.019	99	88592	19.333	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.177	67	34618	19.338	ng/ul	0.00
11) 2-Chlorophenol-d4	7.383	132	86184	19.142	ng/ul	0.00
15) 4-Methylphenol-d8	8.558	113	82008	19.824	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	42540	18.820	ng/ul	0.00
24) 2-Nitrophenol-d4	9.721	143	55961	19.386	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.262	165	116576	19.349	ng/ul	0.00
31) 4-Chloroaniline-d4	10.773	131	128758	18.212	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	434416	19.918	ng/ul	0.00
49) Acenaphthylene-d8	14.175	160	485998	21.027	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	64328	19.556	ng/ul	0.00
60) Fluorene-d10	15.479	176	418827	20.958	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.597	200	94048	17.863	ng/ul	0.00
73) Anthracene-d10	17.330	188	769748	21.297	ng/ul	0.00
81) Pyrene-d10	19.627	212	1016631	22.507	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.316	264	1083008	21.417	ng/ul	0.00
Target Compounds						
71) Pentachlorophenol	16.889	266	32734m	5.107	ng/ul	Qvalue

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

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