

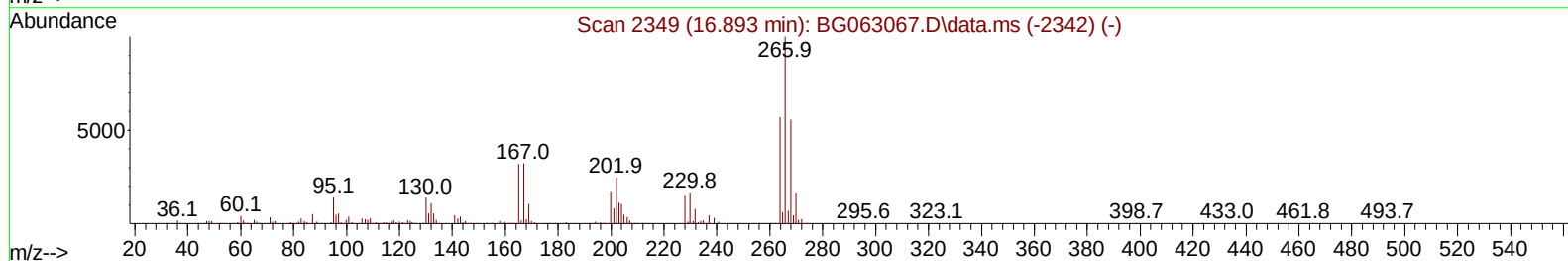
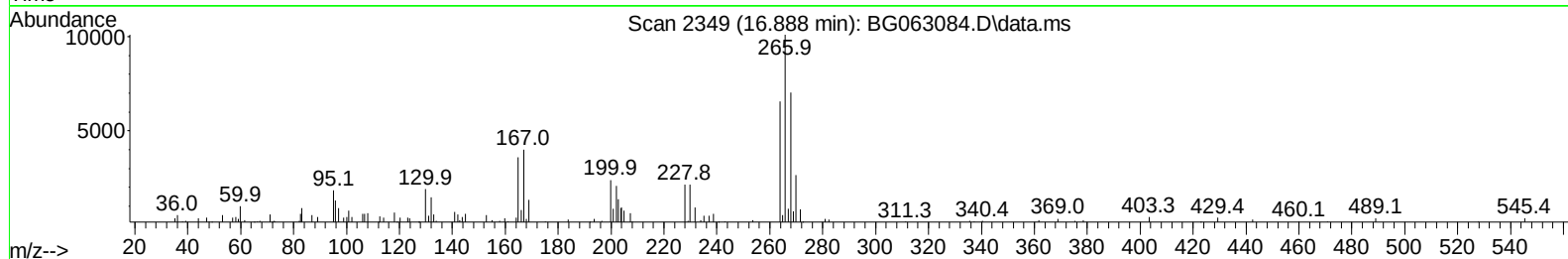
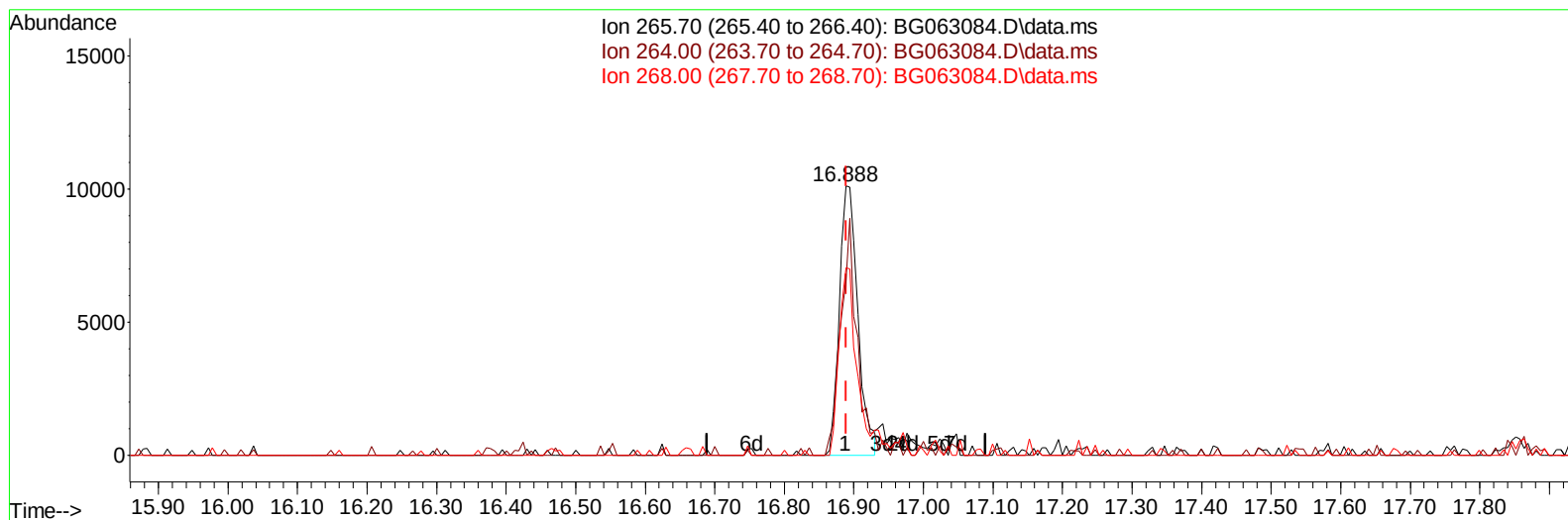
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063084.D
Acq On : 27 Sep 2024 21:46
Operator : RC/JU
Sample : P3910-14
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC34

Manual IntegrationsAPPROVED

Quant Time: Sep 27 22:57:59 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

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



TIC: BG063084.D\data.ms

(71) Pentachlorophenol (C)

16.888min (-0.001) 2.60 ng/uL

response 18757

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	72.10	65.05
268.00	71.60	69.76
0.00	0.00	0.00

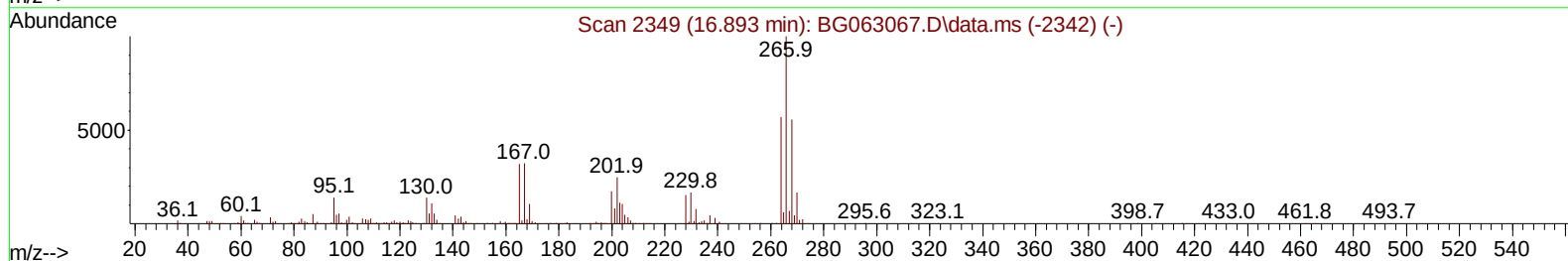
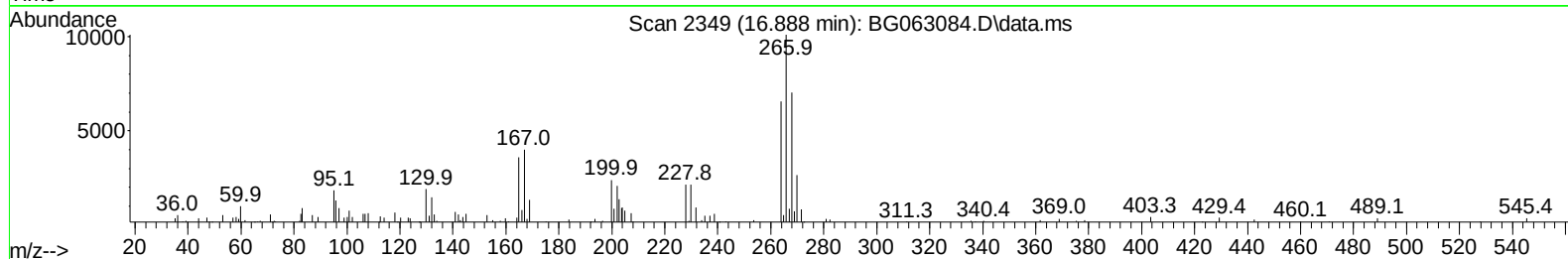
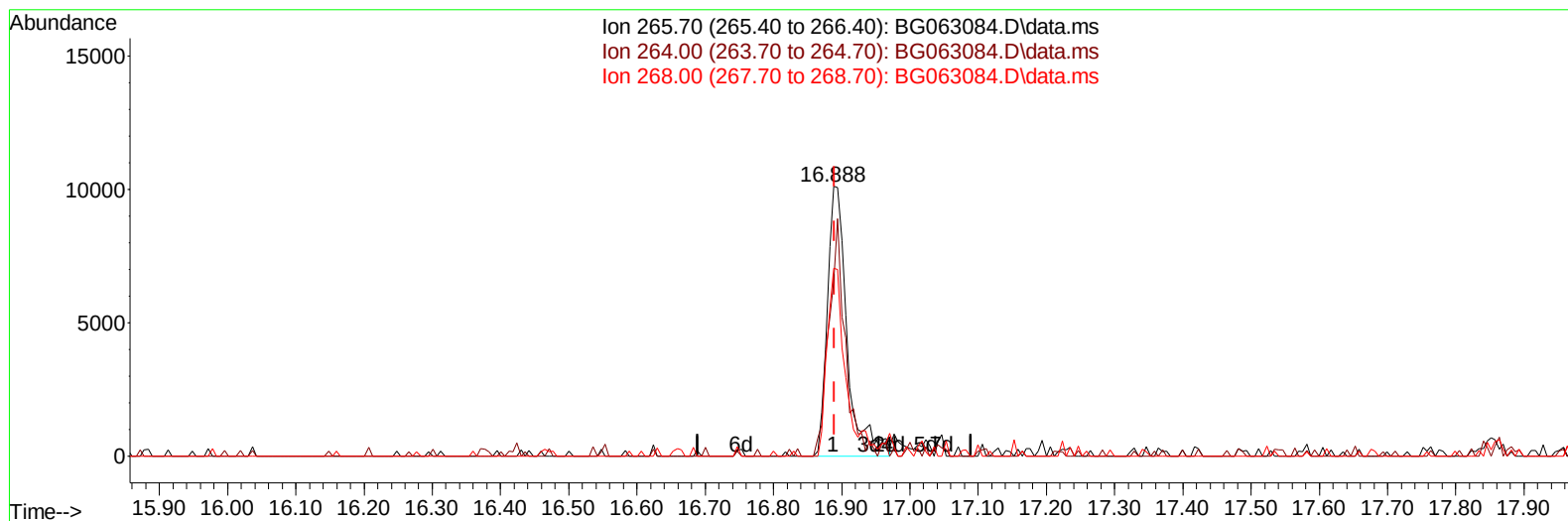
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TIC: BG063084.D\data.ms

(71) Pentachlorophenol (C)

16.888min (-0.001) 2.84 ng/ul m

response 20460

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	72.10	65.05
268.00	71.60	69.76
0.00	0.00	0.00

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 Sample : P3910-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 DDC34

Manual IntegrationsAPPROVED

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

Quant Time: Sep 28 00:41:13 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

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.852	152		73656	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.637	136		316739	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.485	164		307679	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.235	188		847106	20.000	ng/ul	0.00
79) Chrysene-d12	21.489	240		898658	20.000	ng/ul	0.00
88) Perylene-d12	24.520	264		1021324	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.316	96		4752	3.719	ng/uL	0.00
4) Pyridine-d5	3.721	84		67125	21.486	ng/ul	0.00
7) Phenol-d5	7.017	99		105522	23.195	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.182	67		41609	23.411	ng/ul	0.00
11) 2-Chlorophenol-d4	7.382	132		95770	21.425	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113		95343	23.215	ng/ul	0.00
21) Nitrobenzene-d5	8.998	128		47991	20.605	ng/ul	0.00
24) 2-Nitrophenol-d4	9.726	143		67870	22.817	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.261	165		138883	22.371	ng/ul	0.00
31) 4-Chloroaniline-d4	10.772	131		168367	23.112	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166		578878	23.663	ng/ul	0.00
49) Acenaphthylene-d8	14.174	160		604104	23.303	ng/ul	0.00
54) 4-Nitrophenol-d4	14.691	143		60010	16.265	ng/ul	0.00
60) Fluorene-d10	15.478	176		560688	25.014	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.602	200		70874	11.972	ng/ul	0.00
73) Anthracene-d10	17.335	188		1075154	26.455	ng/ul	0.00
81) Pyrene-d10	19.632	212		1381527	27.302	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.315	264		1449455	26.498	ng/ul	0.00
Target Compounds							
71) Pentachlorophenol	16.888	266		20460m	2.839	ng/ul	Qvalue

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

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