

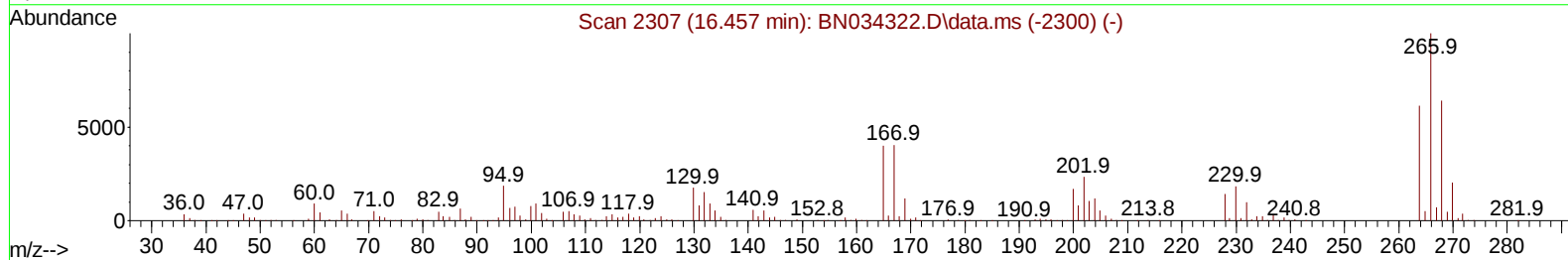
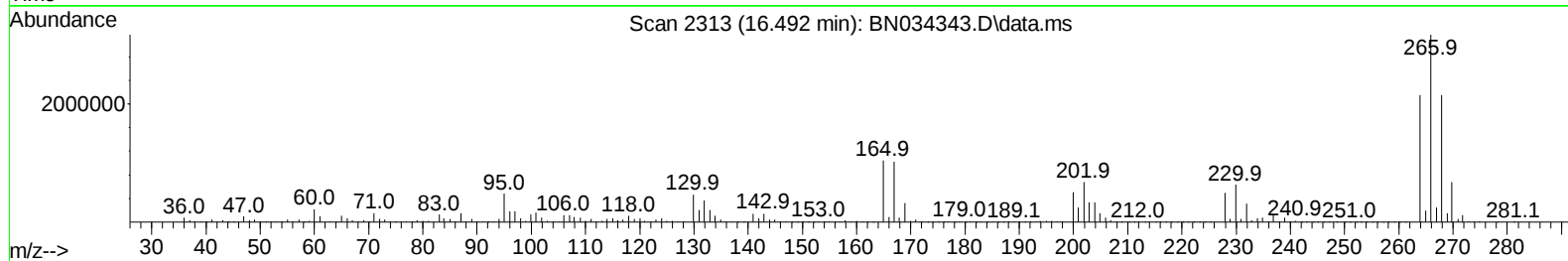
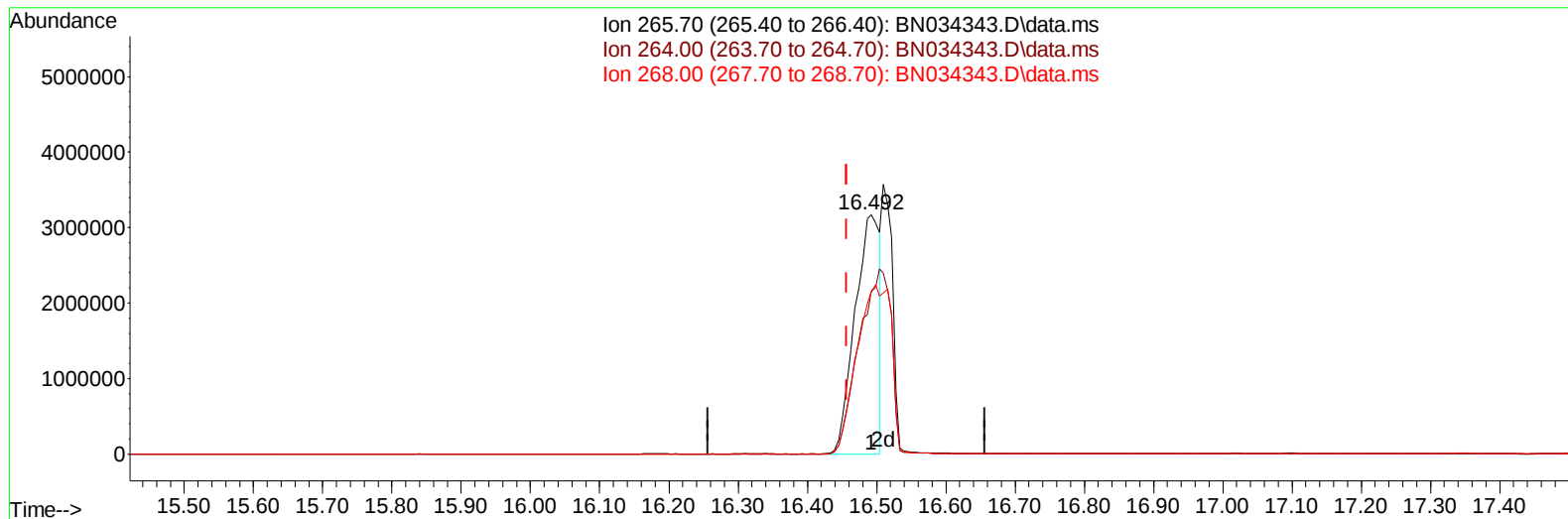
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034343.D
Acq On : 02 Oct 2024 18:55
Operator : RC/JU
Sample : P4016-21
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDC24

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/04/2024
Supervised By :mohammad ahmed 10/05/2024

Quant Time: Oct 02 23:05:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Sep 30 14:14:38 2024
Response via : Initial Calibration



TIC: BN034343.D\data.ms

(71) Pentachlorophenol (C)

16.492min (+ 0.035) 789.33 ng/u1

response 7788565

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.90	67.73
268.00	66.00	67.86
0.00	0.00	0.00

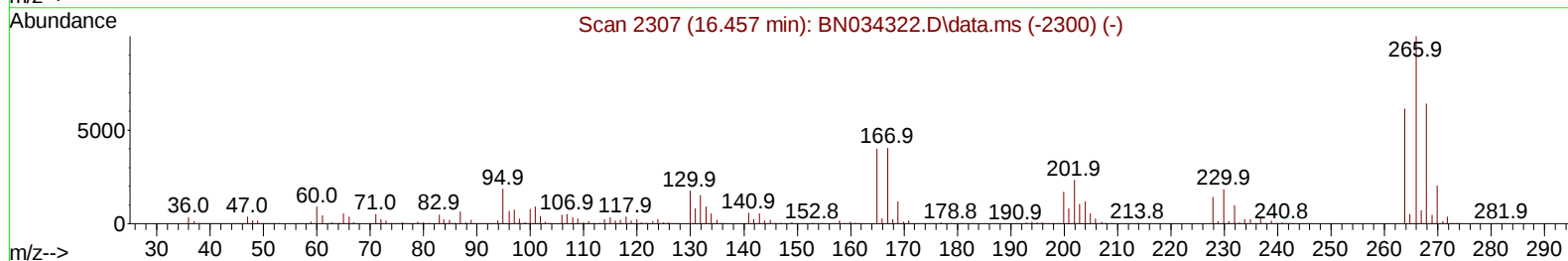
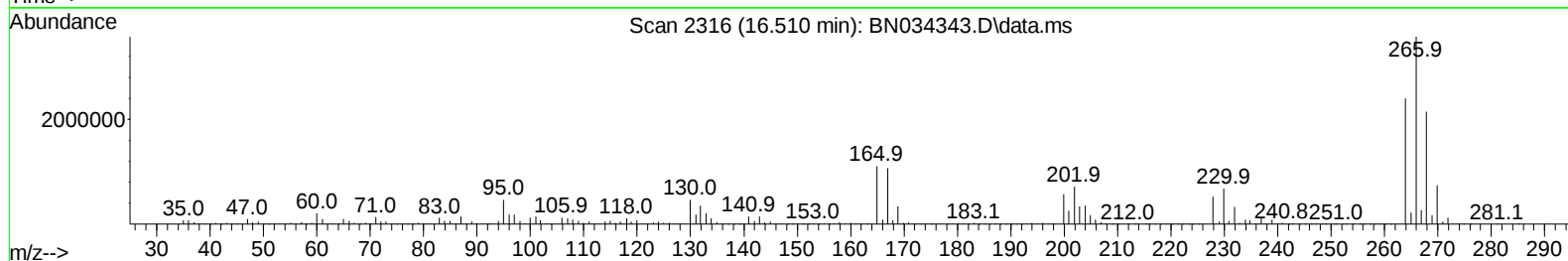
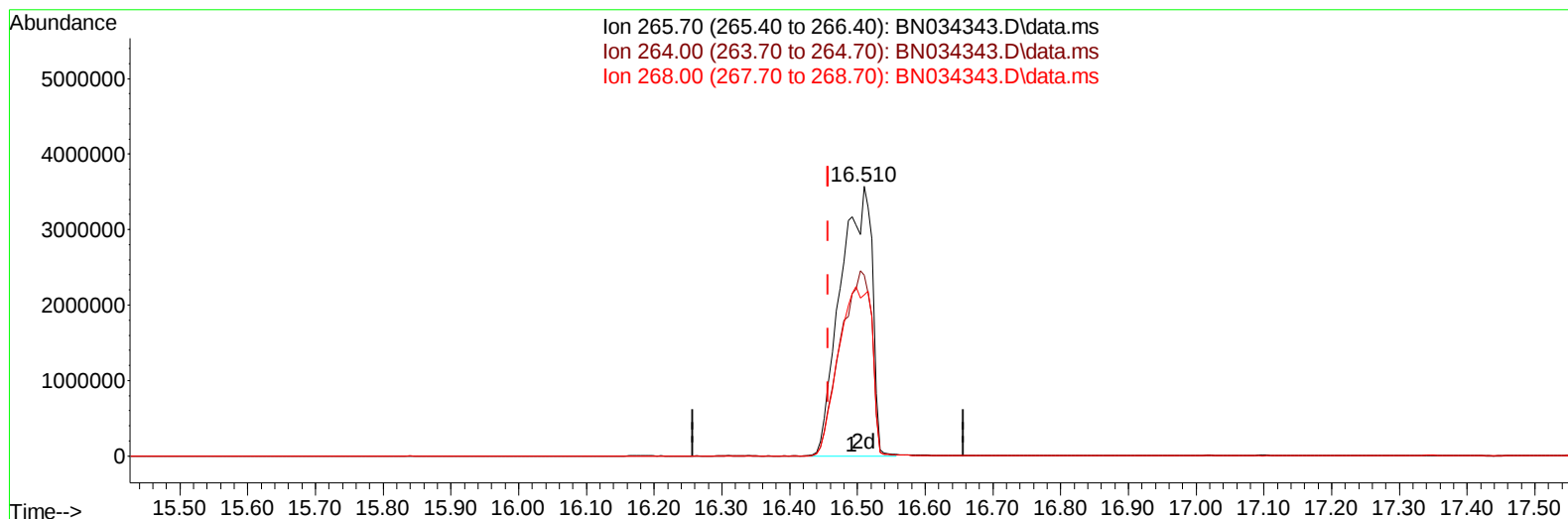
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TIC: BN034343.D\data.ms

(71) Pentachlorophenol (C)

16.510min (+ 0.053) 1174.96 ng/ul m

response 11593774

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.90	67.08
268.00	66.00	59.85
0.00	0.00	0.00

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 QLast Update : Mon Sep 30 14:14:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.393	152	228121	20.000	ng/ul	0.00
20) Naphthalene-d8	10.146	136	857972	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.045	164	621222	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.810	188	1257009	20.000	ng/ul	0.01
79) Chrysene-d12	21.033	240	1240149	20.000	ng/ul	0.00
88) Perylene-d12	23.751	264	1504382	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.999	96	18582	3.324	ng/uL	0.00
4) Pyridine-d5	3.393	84	230135	13.850	ng/ul	0.00
7) Phenol-d5	6.593	99	383193	18.927	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.740	67	253603	20.189	ng/ul	0.00
11) 2-Chlorophenol-d4	6.934	132	330384	20.731	ng/ul	0.00
15) 4-Methylphenol-d8	8.110	113	302550	18.162	ng/ul	0.00
21) Nitrobenzene-d5	8.534	128	162351	23.741	ng/ul	0.00
24) 2-Nitrophenol-d4	9.246	143	175933	25.018	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.787	165	307250	23.200	ng/ul	0.01
31) 4-Chloroaniline-d4	10.298	131	210977	10.523	ng/ul	0.00
46) Dimethylphthalate-d6	13.469	166	897016	19.321	ng/ul	0.00
49) Acenaphthylene-d8	13.734	160	1034930	19.815	ng/ul	0.00
54) 4-Nitrophenol-d4	14.286	143	164524	17.268	ng/ul	0.01
60) Fluorene-d10	15.051	176	810572	20.601	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.181	200	115872	15.460	ng/ul	0.00
73) Anthracene-d10	16.910	188	1200345	20.311	ng/ul	0.01
81) Pyrene-d10	19.210	212	1377762	19.939	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	1552256	20.087	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.198	128	1089374	23.213	ng/ul	97
36) 2-Methylnaphthalene	11.828	142	2111728	65.325	ng/ul	99
37) 1-Methylnaphthalene	12.051	142	1258880	38.472	ng/ul	97
41) 2,4,6-Trichlorophenol	12.457	196	11946	1.005	ng/ul	89
42) 2,4,5-Trichlorophenol	12.534	196	176128	13.487	ng/ul	99
43) 1,1'-Biphenyl	12.863	154	259926	5.550	ng/ul	91
52) Acenaphthene	14.110	153	78181	1.916	ng/ul	90
58) 2,3,4,6-Tetrachlorophenol	14.692	232	1467475	135.414	ng/ul	90
61) Fluorene	15.104	166	188591	4.132	ng/ul#	87
71) Pentachlorophenol	16.510	266	11593774m	1174.962	ng/ul	
72) Phenanthrene	16.857	178	786520	11.348	ng/ul#	96
86) Bis(2-ethylhexyl)phtha...	20.980	149	68287	1.198	ng/ul#	89

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

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