Quantitation Report (Qedit)

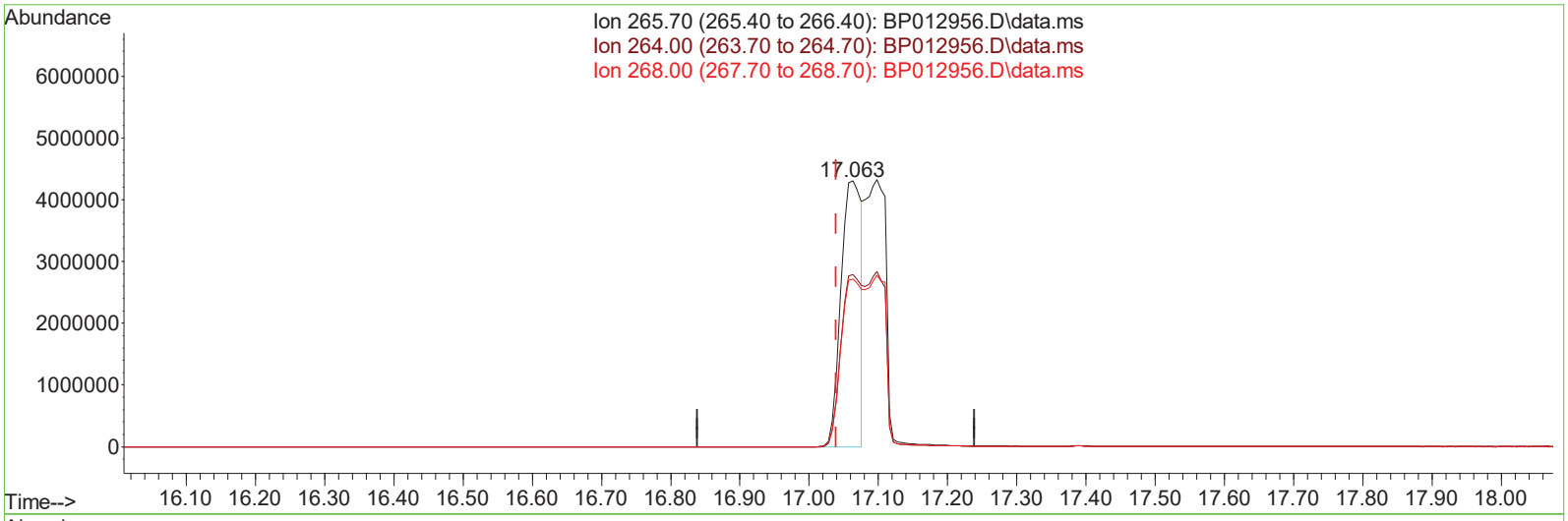
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012953.D
Acq On : 05 Dec 2022 15:31
Operator : CG/JU
Sample : N5738-15DL2 30X
Misc :
ALS Vial : 9 Sample Multiplier: 1

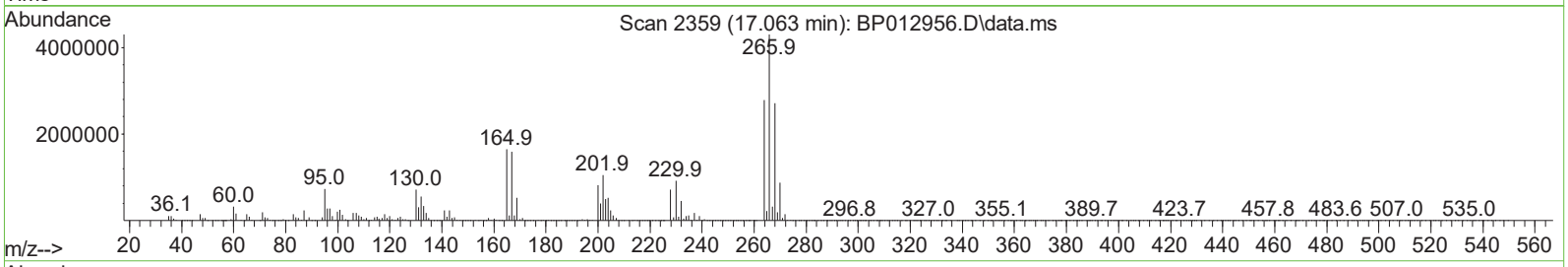
Quant Time: Dec 05 21:51:45 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 01 23:46:27 2022
Response via : Initial Calibration

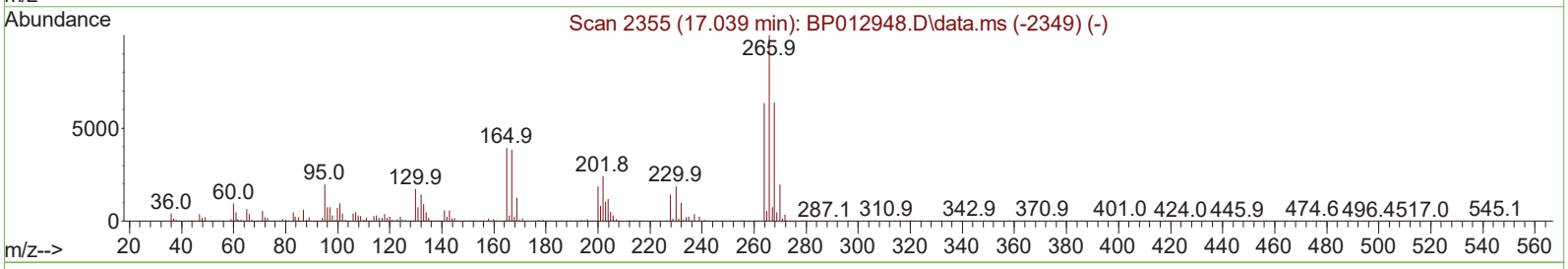
Instrument :
BNA_P
ClientSampleId :
GBNJ4DL2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2022
Supervised By :mohammad ahmed 12/08/2022







TIC: BP012956.D\data.ms

(71) Pentachlorophenol (C)

17.063min (+ 0.024) 284.06 ng/ul

response 8664673

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.10	64.71
268.00	64.00	63.06
0.00	0.00	0.00

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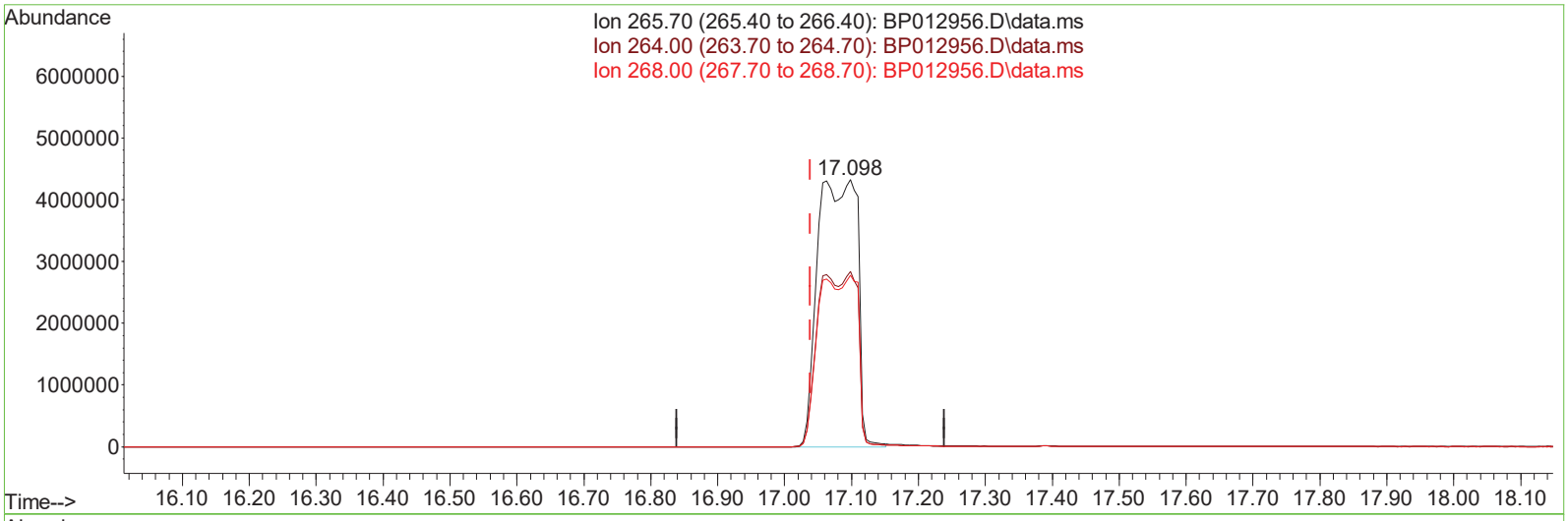

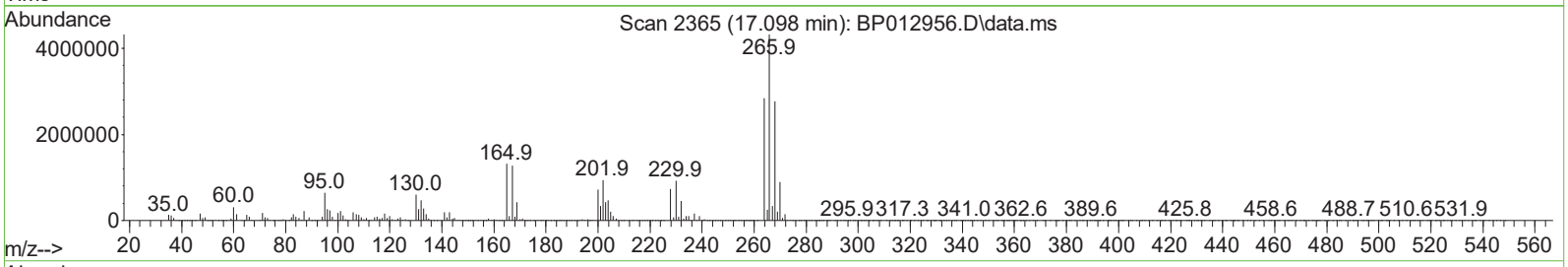

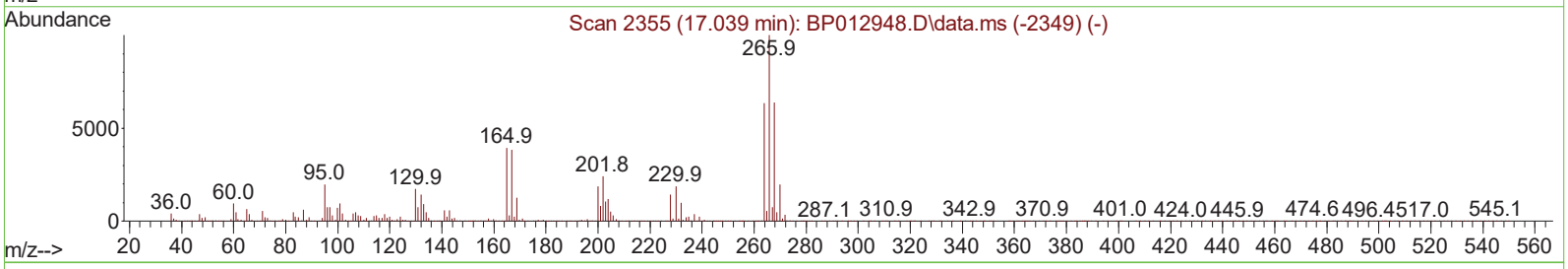

TIC: BP012956.D\data.ms

(71) Pentachlorophenol (C)

17.098min (+ 0.059) 581.95 ng/ul m

response 17751412

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.10	65.60
268.00	64.00	64.19
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012953.D
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Sample : N5738-15DL2 30X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 05 21:51:45 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
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Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2022
Supervised By :mohammad ahmed 12/08/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	568258	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2581546	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1772858	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.416	188	3808913	20.000	ng/ul	0.02
79) Chrysene-d12	21.492	240	3130343	20.000	ng/ul	0.01
88) Perylene-d12	24.010	264	2680620	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	913	0.071	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.158	99	10920	0.208	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	9082	0.306	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	10279	0.267	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	8870	0.206	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	4854	0.264	ng/ul	0.00
24) 2-Nitrophenol-d4	9.905	143	5045	0.240	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	9888	0.235	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	14.057	166	48604	0.380	ng/ul	0.01
49) Acenaphthylene-d8	14.345	160	44826	0.318	ng/ul	0.01
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.645	176	43702	0.389	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.751	200	3493	0.156	ng/ul	0.00
73) Anthracene-d10	17.510	188	67670	0.402	ng/ul	0.01
81) Pyrene-d10	19.733	212	66098	0.430	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.839	264	40573	0.310	ng/ul	0.01
Target Compounds						Qvalue
58) 2,3,4,6-Tetrachlorophenol	15.275	232	6015976	161.026	ng/ul	95
71) Pentachlorophenol	17.098	266	17751412m	581.948	ng/ul	

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

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