Quantitation Report (Qedit)

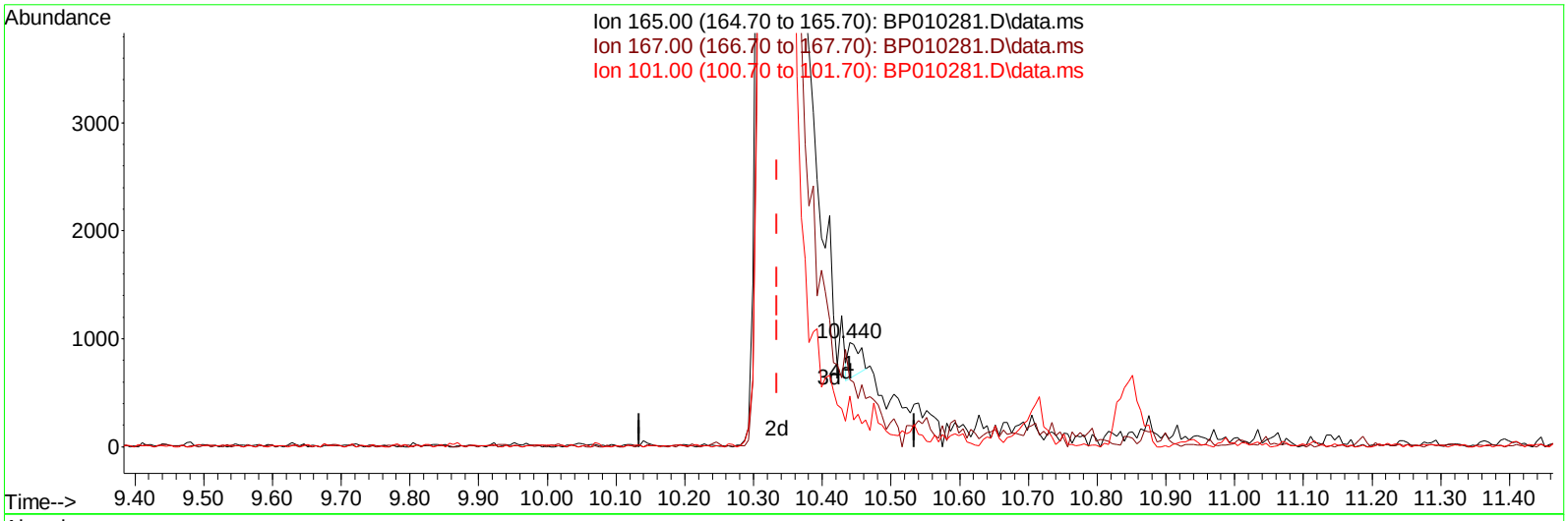
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010281.D
Acq On : 11 May 2022 10:21
Operator : CG/JU
Sample : PB144607BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

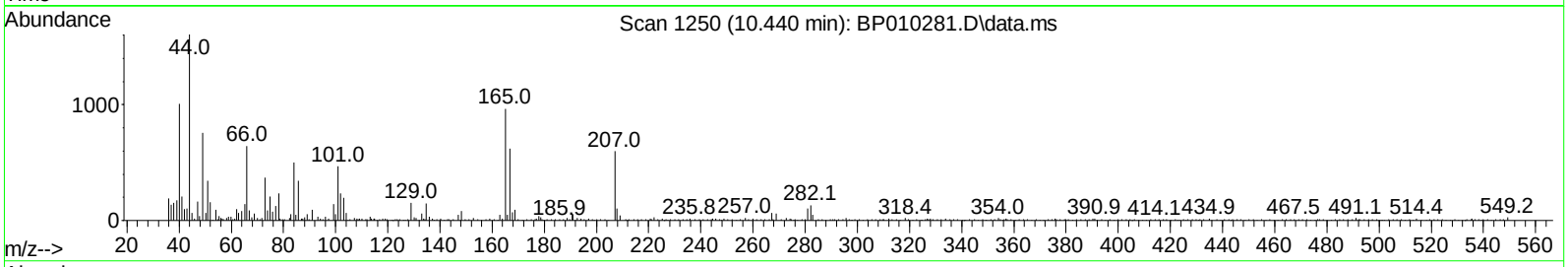
Quant Time: May 12 01:11:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

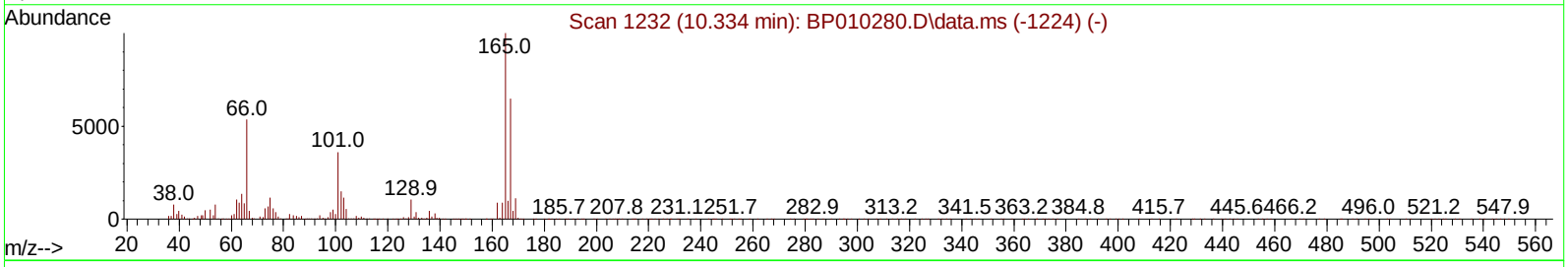
Instrument :
BNA_P
ClientSampleId :
SBLK607

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022







TIC: BP010281.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.440min (+ 0.106) 0.04 ng/ul

response 382

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	64.25
101.00	49.80	48.50
0.00	0.00	0.00

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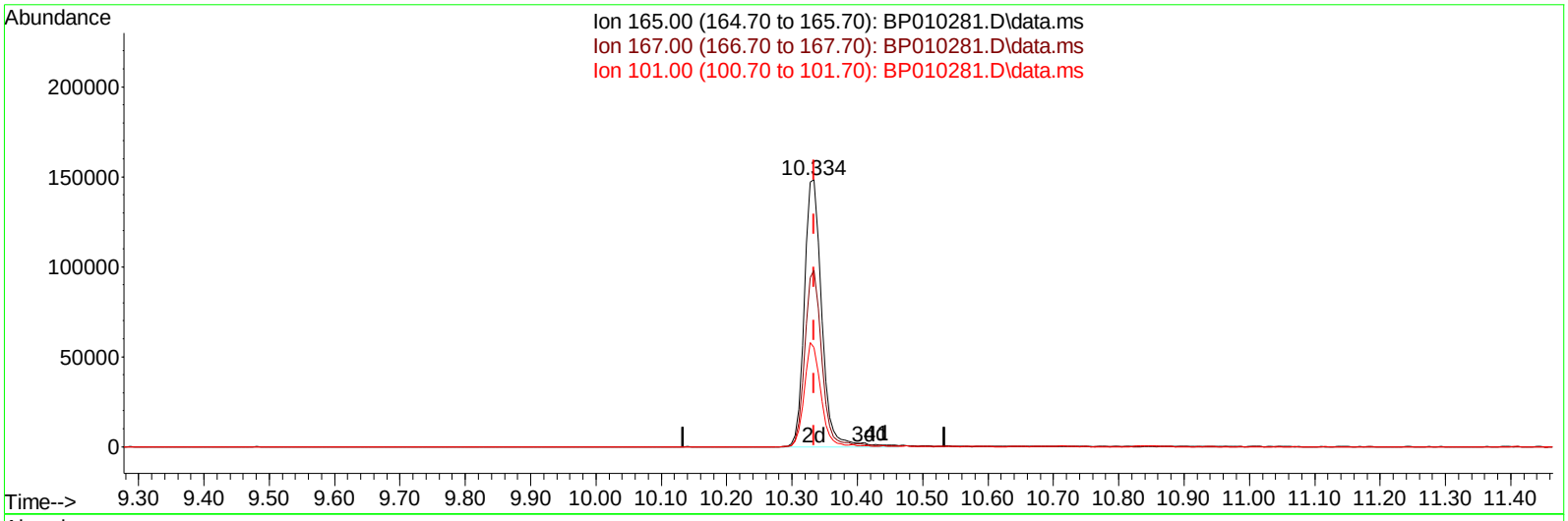

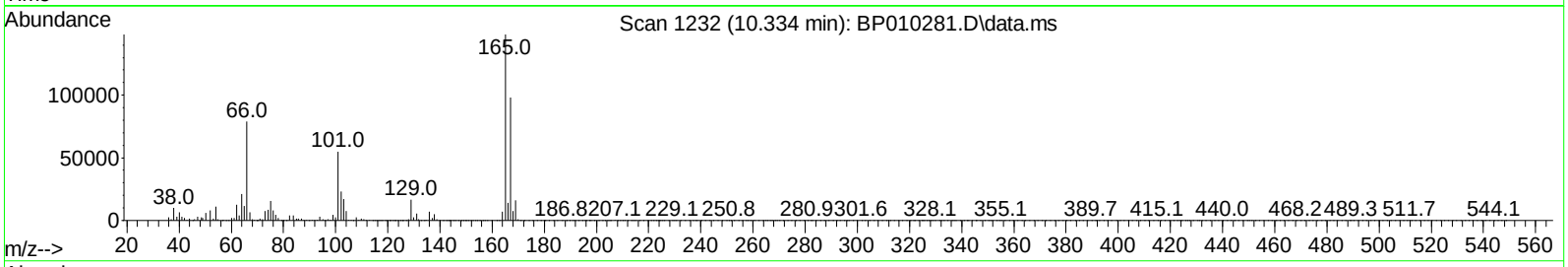

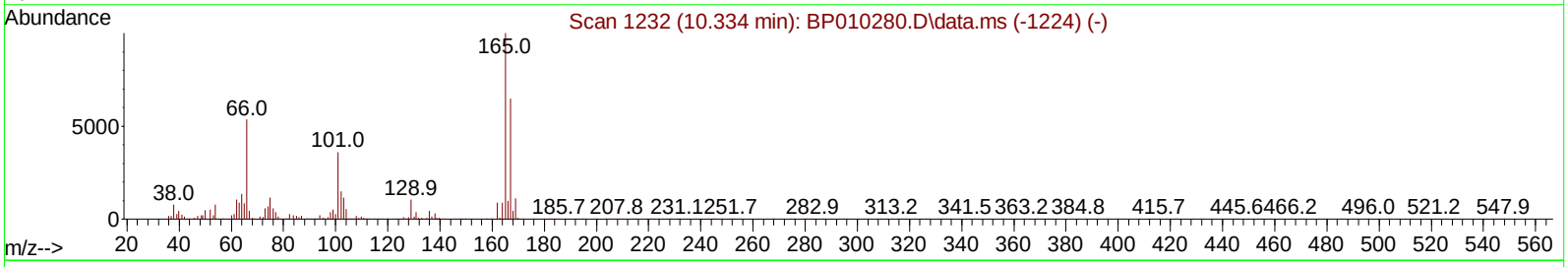

TIC: BP010281.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.334min (+ 0.000) 27.39 ng/ul m

response 272042

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	66.15
101.00	49.80	37.07#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
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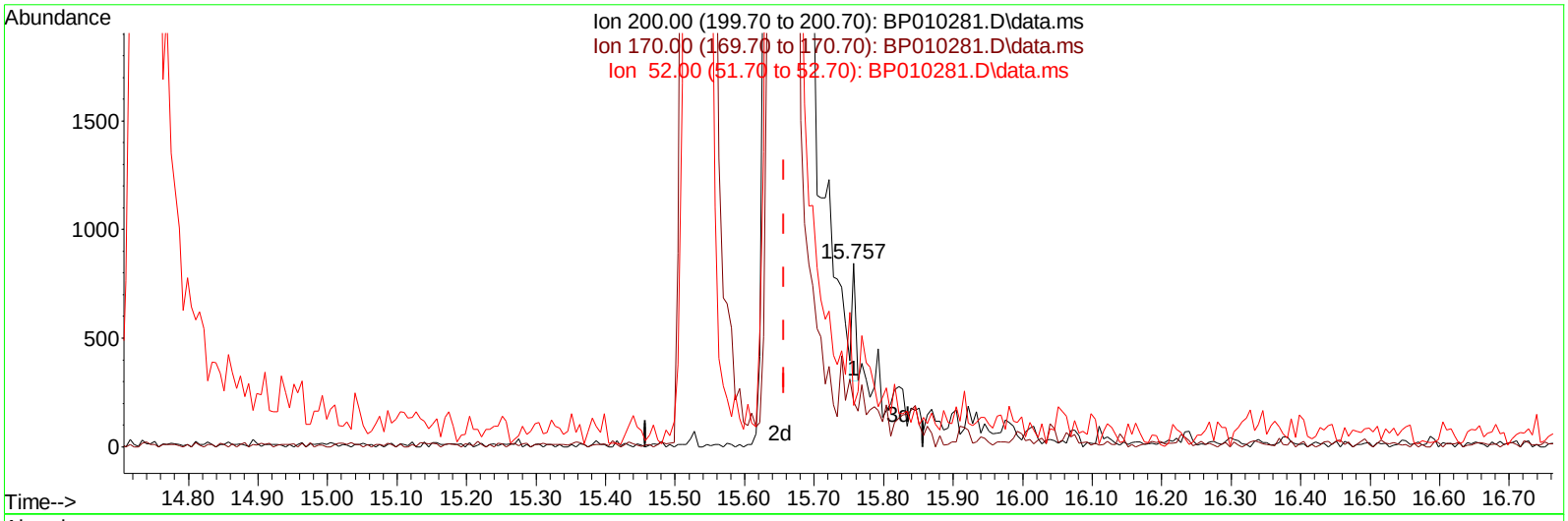

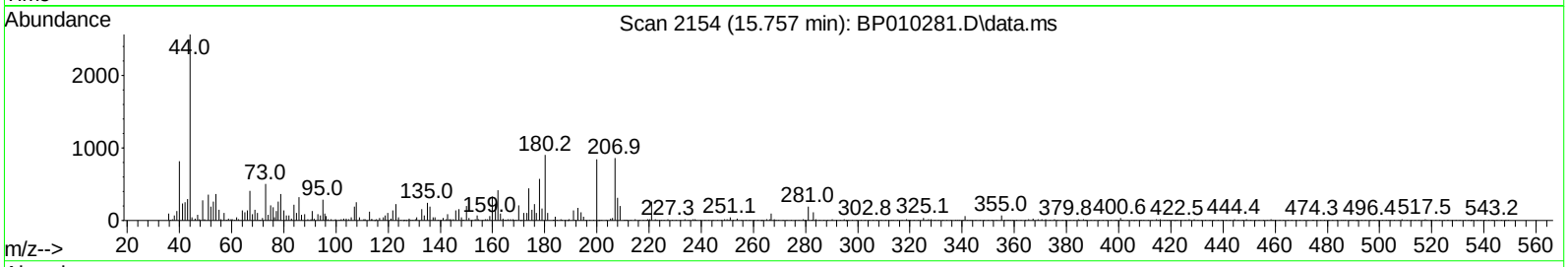

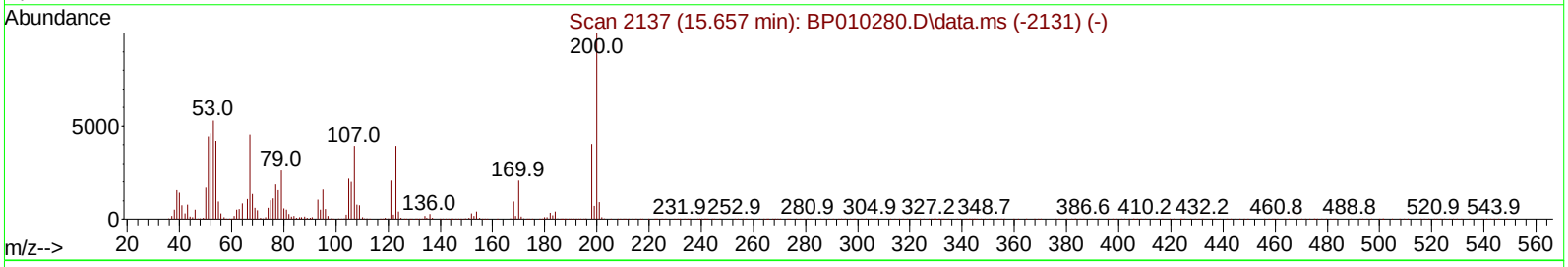

TIC: BP010281.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.757min (+ 0.100) 0.03 ng/ul

response 191

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	24.62
52.00	21.20	22.37
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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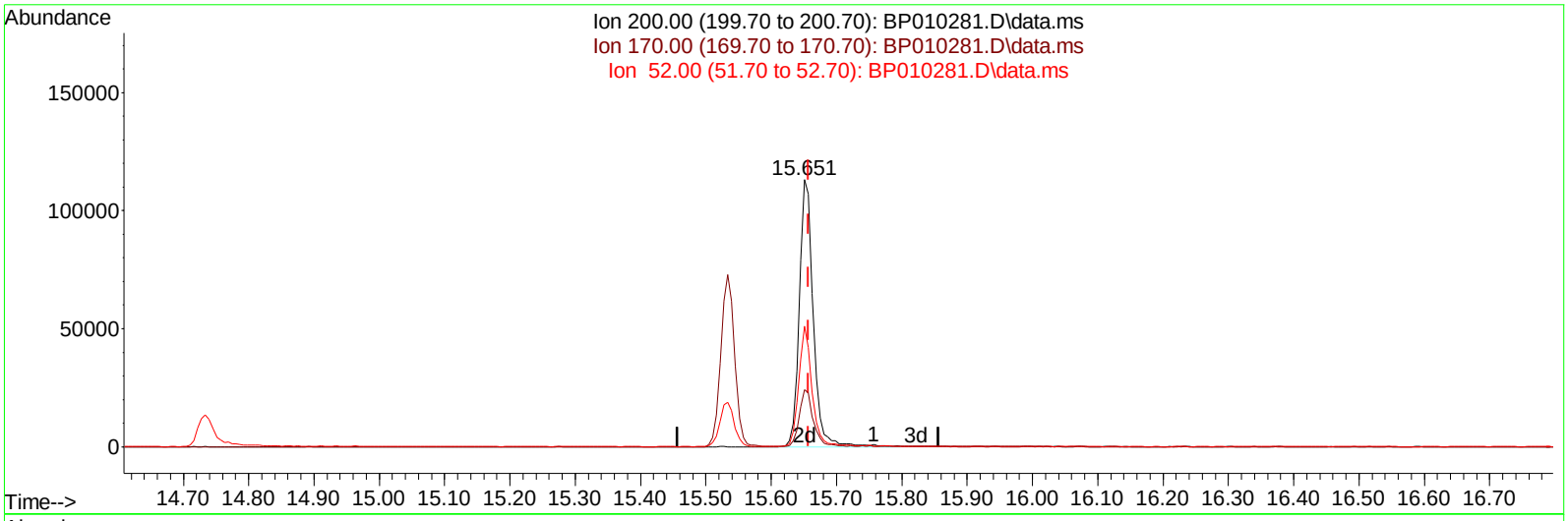

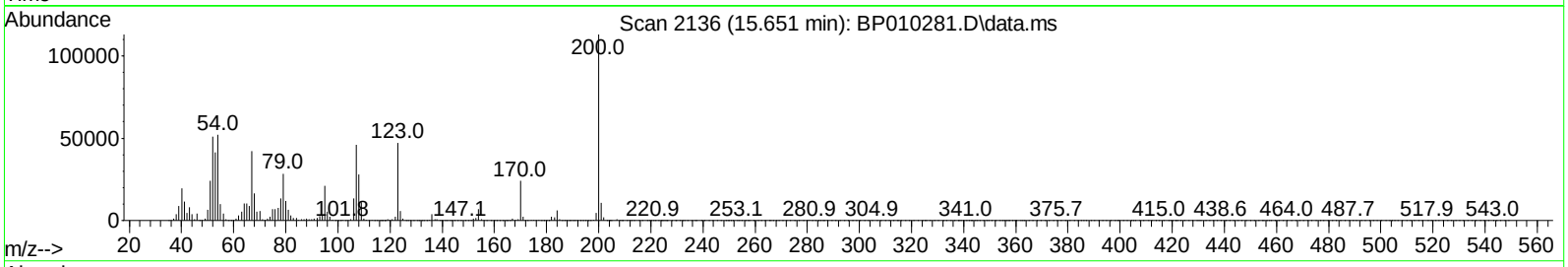

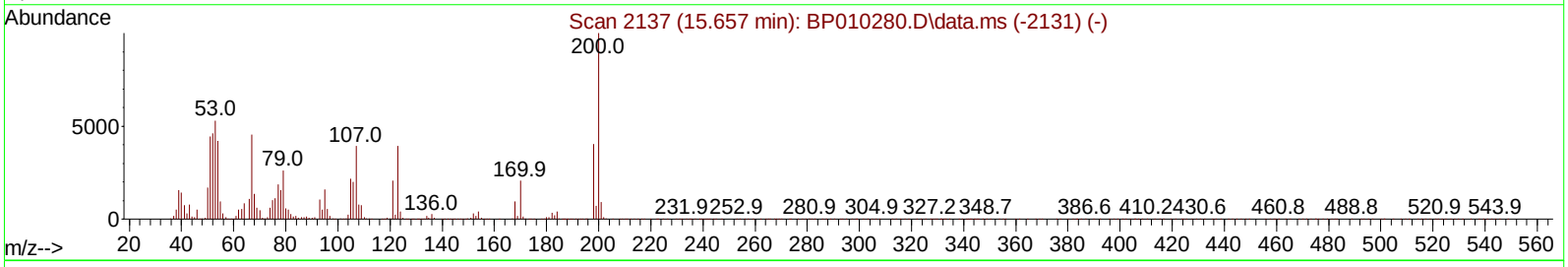

TIC: BP010281.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.651min (-0.006) 25.93 ng/ul m

response 167839

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	21.29
52.00	21.20	45.16#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 05/17/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	133974	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	596878	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	424857	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	945391	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.374	240	930737	20.000	ng/ul	# 0.00
88) Perylene-d12	23.810	264	877948	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	20650	6.548	ng/uL	0.00
4) Pyridine-d5	3.770	84	274032	31.225	ng/ul	0.00
7) Phenol-d5	7.069	99	340888	27.395	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	245376	31.876	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	278675	30.409	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	294449	28.783	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	147479	30.816	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	162344	29.751	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	272042m	27.391	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	407799	27.813	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	1097635	32.429	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	1173614	29.083	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	155068	23.885	ng/ul	0.00
60) Fluorene-d10	15.534	176	908721	31.806	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	167839m	25.925	ng/ul	0.00
73) Anthracene-d10	17.392	188	1462369	31.587	ng/ul	0.00
81) Pyrene-d10	19.622	212	1784046	34.687	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.657	264	1551056	33.127	ng/ul	0.00
Target Compounds						Qvalue

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

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