

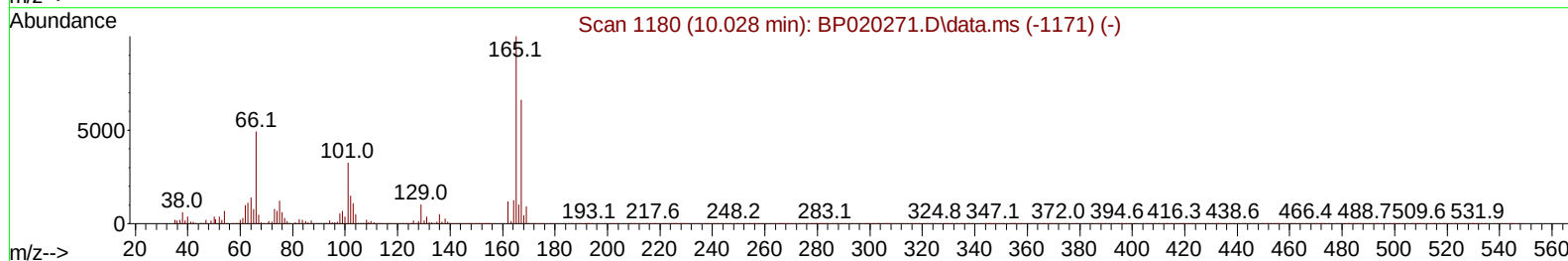
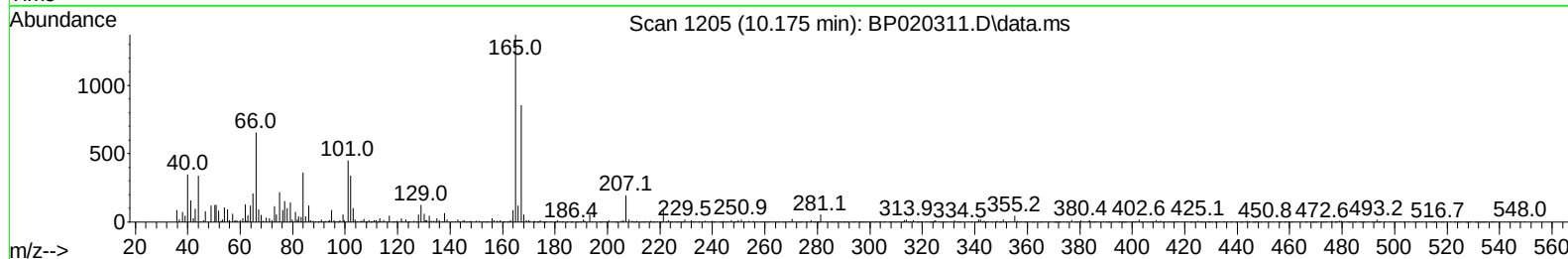
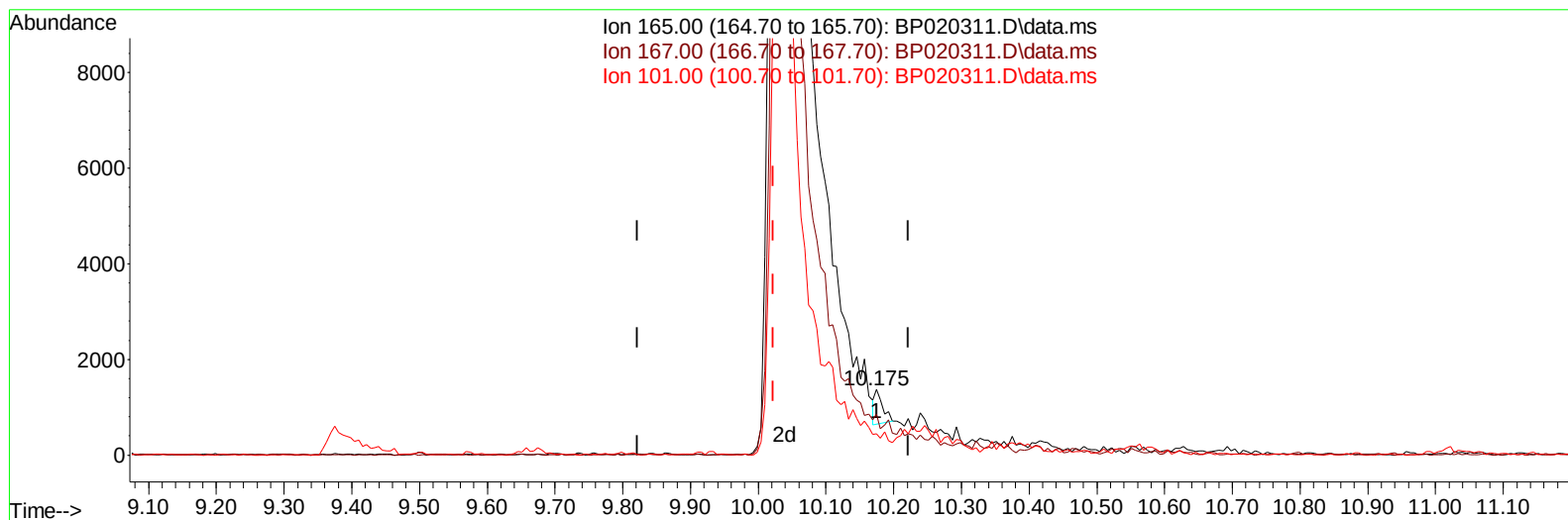
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020311.D
Acq On : 11 May 2024 06:22
Operator : MA/JU
Sample : P2370-14
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Q6

Manual IntegrationsAPPROVED

Quant Time: May 11 07:02:42 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 08 22:57:55 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/13/2024
Supervised By :mohammad ahmed 05/14/2024



TIC: BP020311.D\data.ms

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

10.175min (+ 0.152) 0.13 ng/ul

response 578

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	63.00	62.26
101.00	40.90	32.77
0.00	0.00	0.00

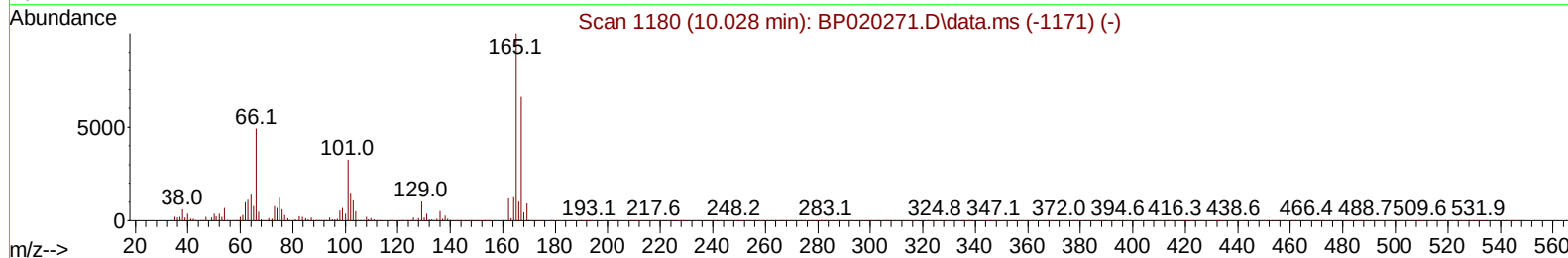
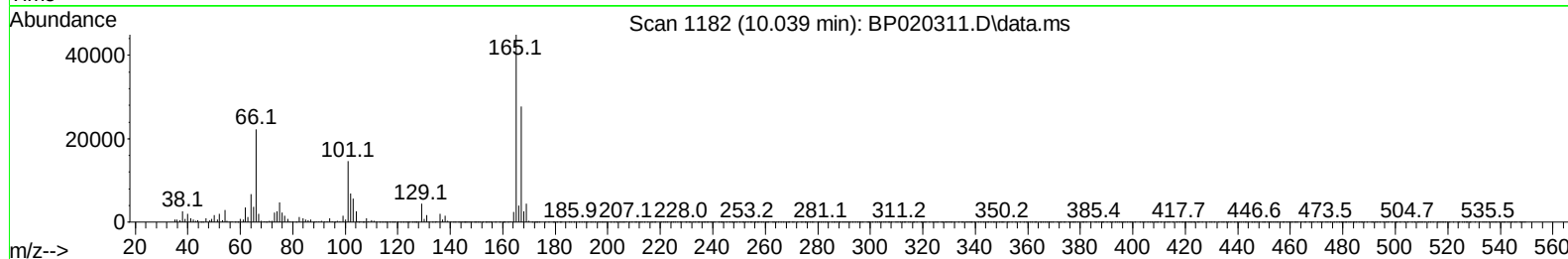
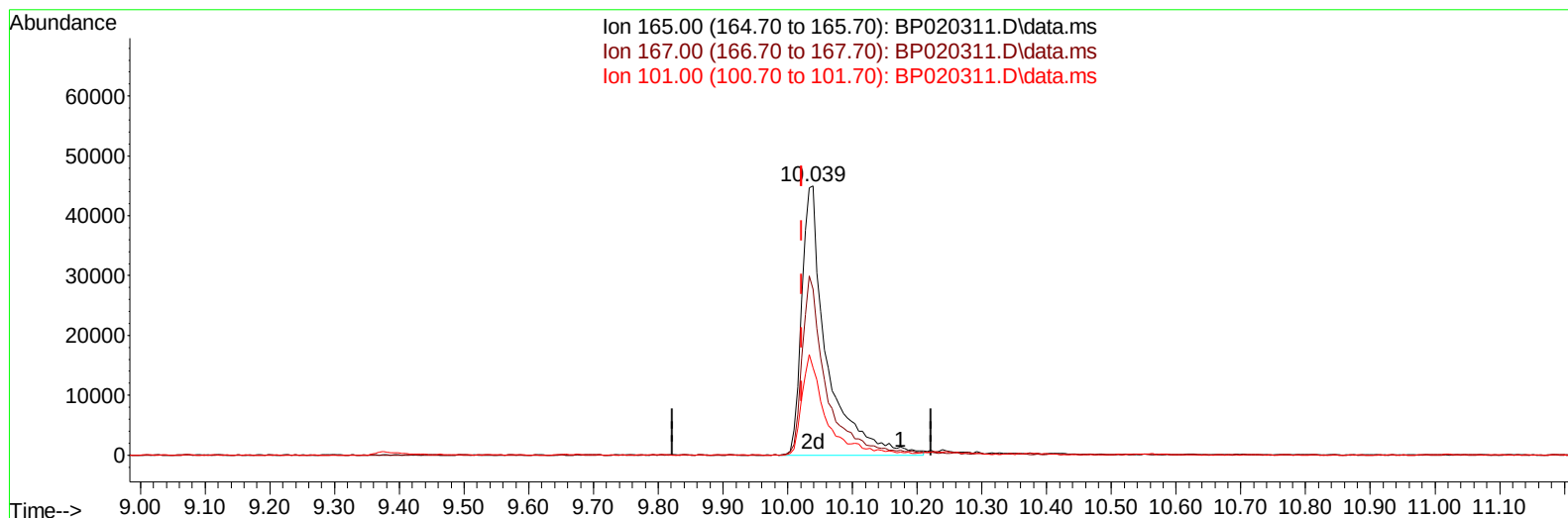
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(28) 2,4-Dichlorophenol-d3 (S)

10.039min (+ 0.017) 27.81 ng/ul m

response 120456

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	63.00	61.73
101.00	40.90	32.46#
0.00	0.00	0.00

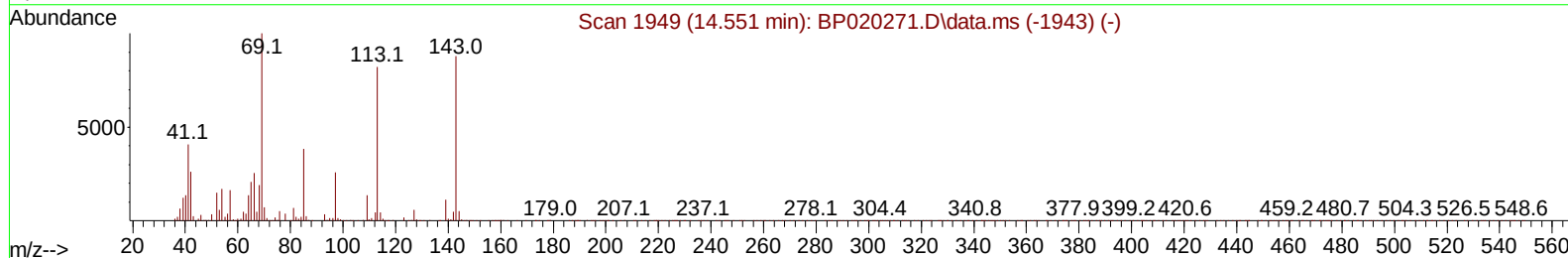
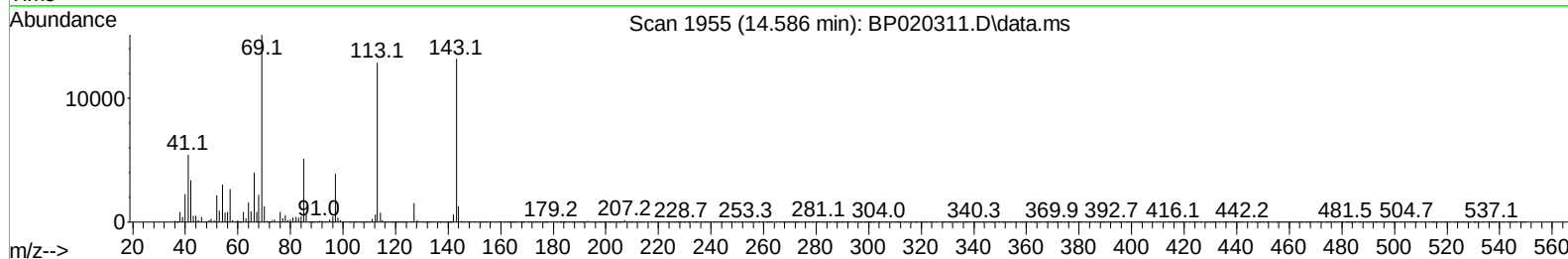
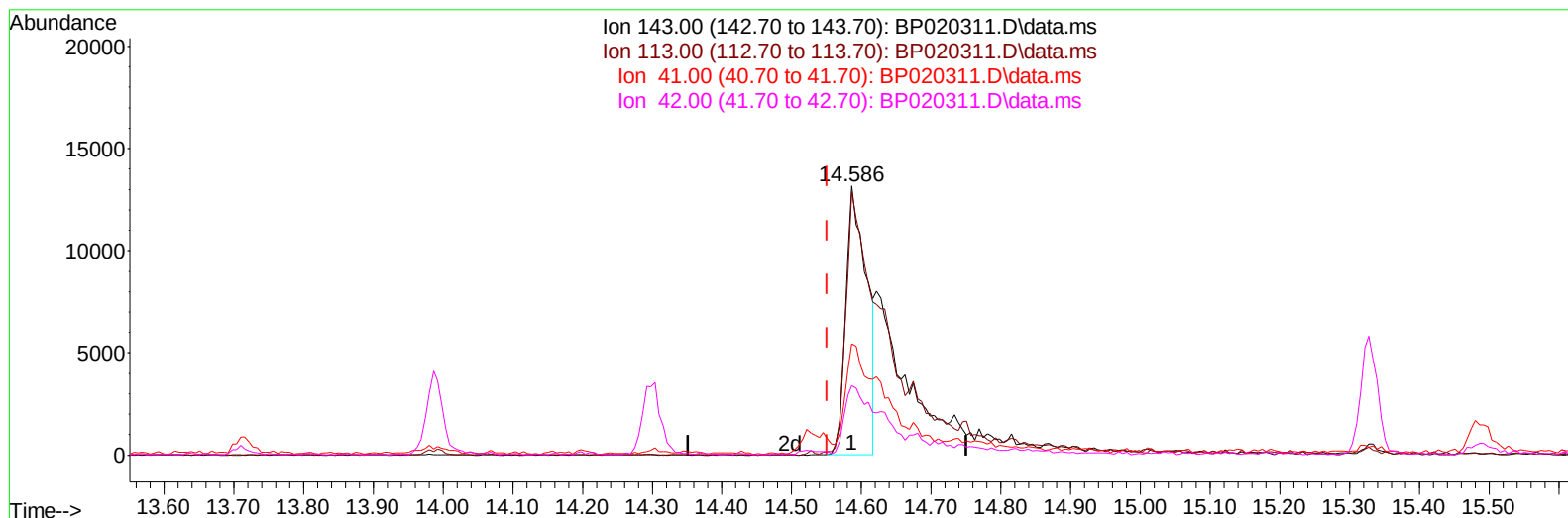
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TIC: BP020311.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.586min (+ 0.035) 12.55 ng/ul

response 27089

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.97
41.00	50.30	41.46
42.00	32.50	25.75#

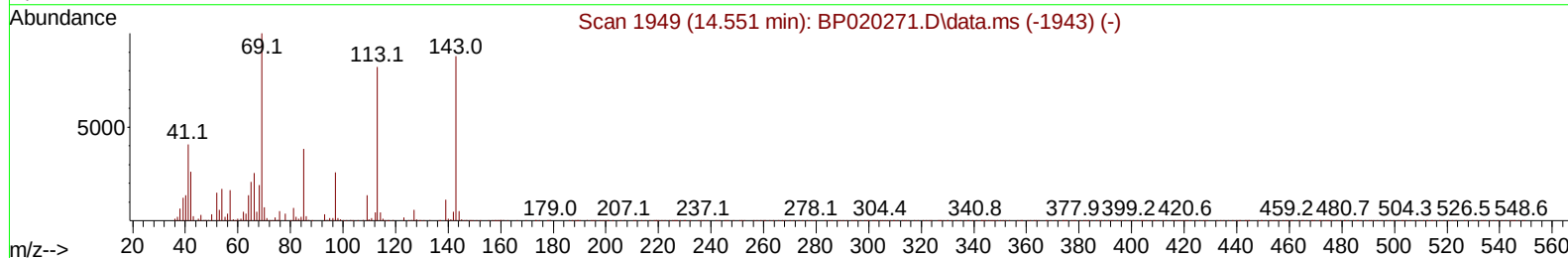
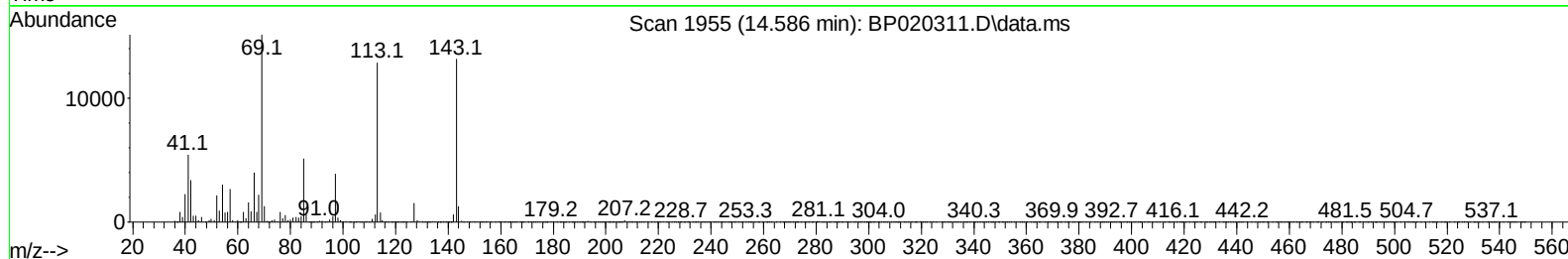
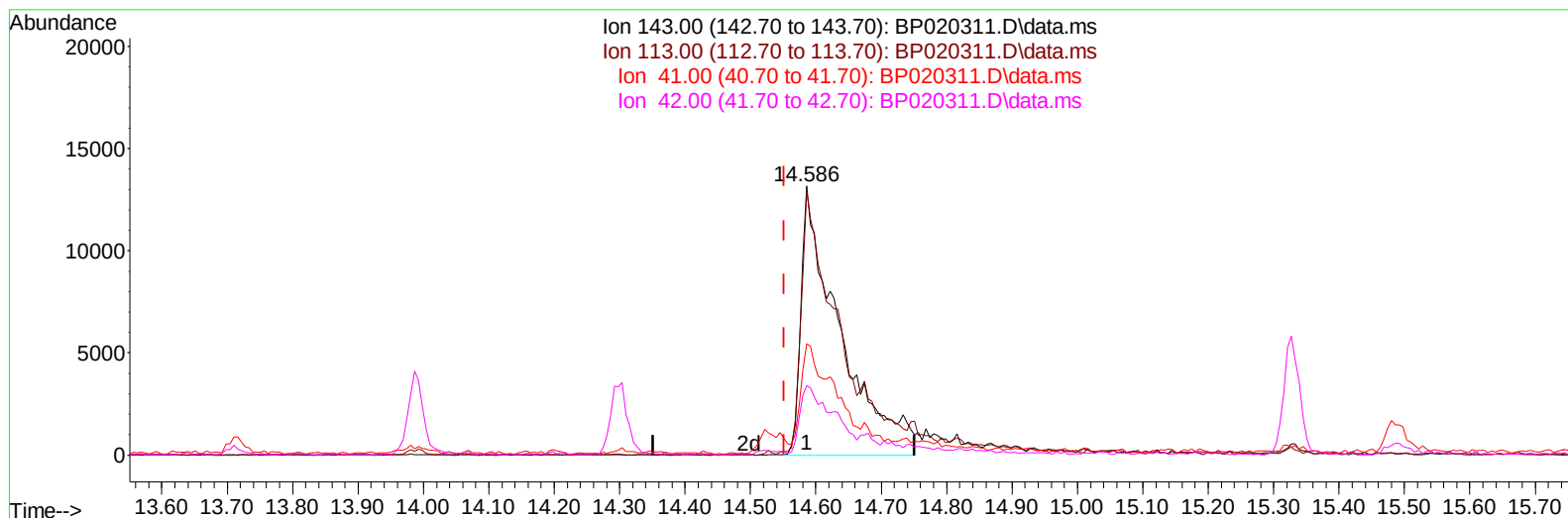
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Instrument :
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TIC: BP020311.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.586min (+ 0.035) 24.92 ng/ul m

response 53787

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.97
41.00	50.30	41.46
42.00	32.50	25.75#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/13/2024
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Quant Time: May 11 07:05:34 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.616	152		66197	20.000	ng/ul	0.00
20) Naphthalene-d8	10.386	136		276803	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164		192305	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.133	188		432714	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240		442744	20.000	ng/ul	0.00
88) Perylene-d12	25.003	264		501334	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.216	96		7492	4.780	ng/uL	0.00
4) Pyridine-d5	3.622	84		96938	21.285	ng/ul	0.00
7) Phenol-d5	6.810	99		141982	24.975	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67		84740	24.538	ng/ul	0.00
11) 2-Chlorophenol-d4	7.151	132		111803	27.510	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113		111958	24.359	ng/ul	0.01
21) Nitrobenzene-d5	8.781	128		56021	27.027	ng/ul	0.00
24) 2-Nitrophenol-d4	9.492	143		66927	28.196	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165		120456m	27.809	ng/ul	0.02
31) 4-Chloroaniline-d4	10.557	131		160774	26.520	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166		420828	29.966	ng/ul	0.00
49) Acenaphthylene-d8	13.986	160		461592	29.755	ng/ul	0.00
54) 4-Nitrophenol-d4	14.586	143		53787m	24.920	ng/ul	0.03
60) Fluorene-d10	15.327	176		362816	30.975	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200		44329	16.144	ng/ul	0.00
73) Anthracene-d10	17.245	188		590356	31.537	ng/ul	0.00
81) Pyrene-d10	19.639	212		732961	28.321	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.768	264		769152	31.731	ng/ul	0.00

Target Compounds Qvalue

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

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