

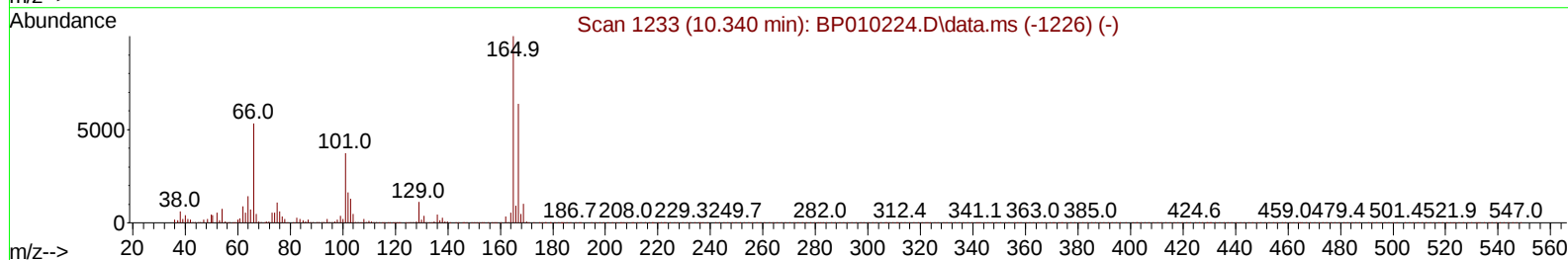
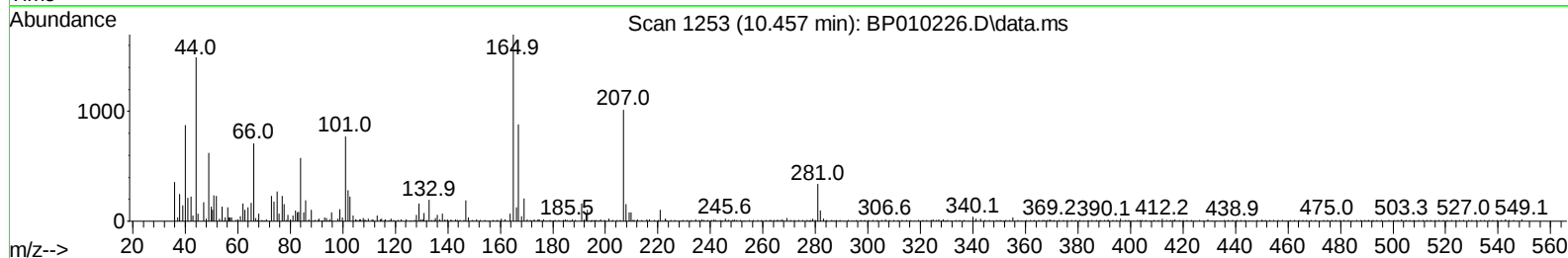
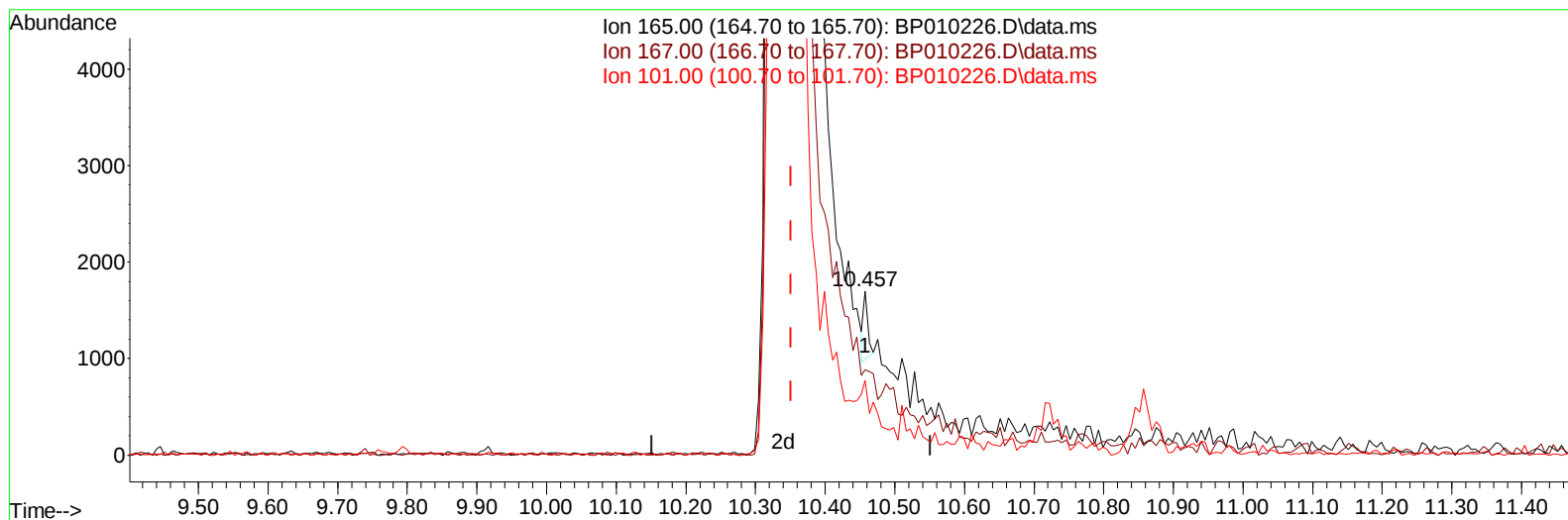
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050422\
Data File : BP010226.D
Acq On : 04 May 2022 16:52
Operator : CG/JU
Sample : PB144568BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK568

Manual IntegrationsAPPROVED

Quant Time: May 05 02:58:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :mohammad ahmed 05/09/2022



TIC: BP010226.D\data.ms

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

10.457min (+ 0.106) 0.03 ng/u1

response 316

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	51.85
101.00	49.80	45.44
0.00	0.00	0.00

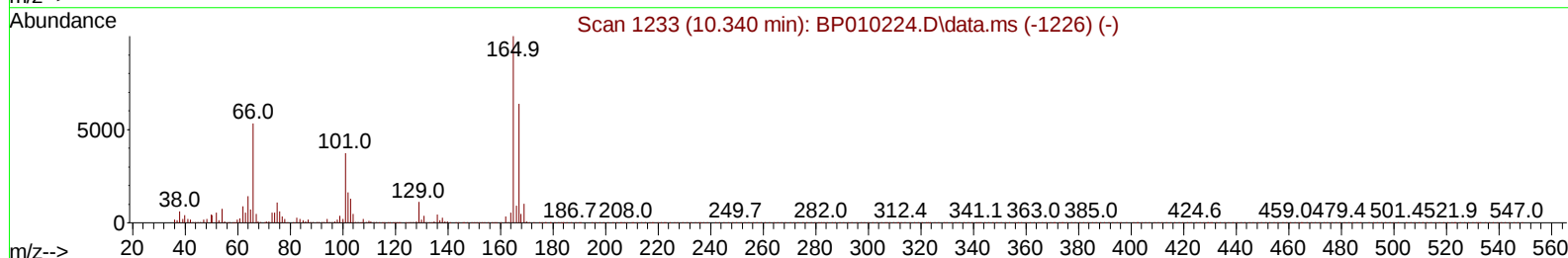
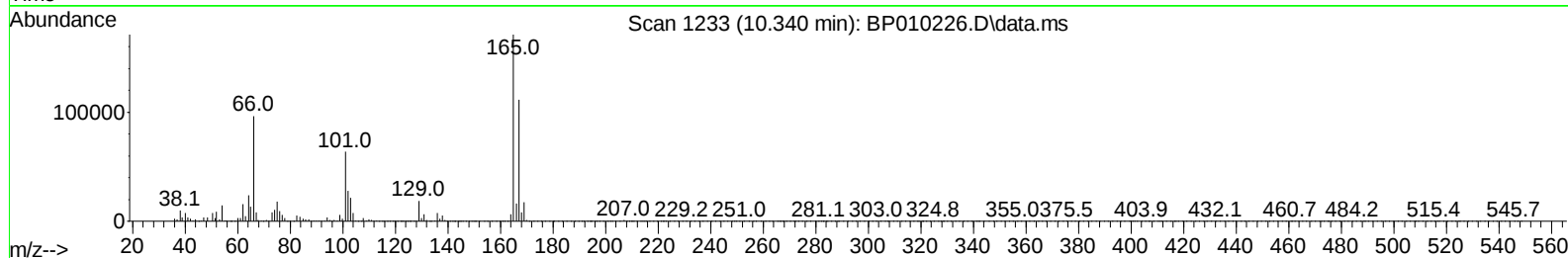
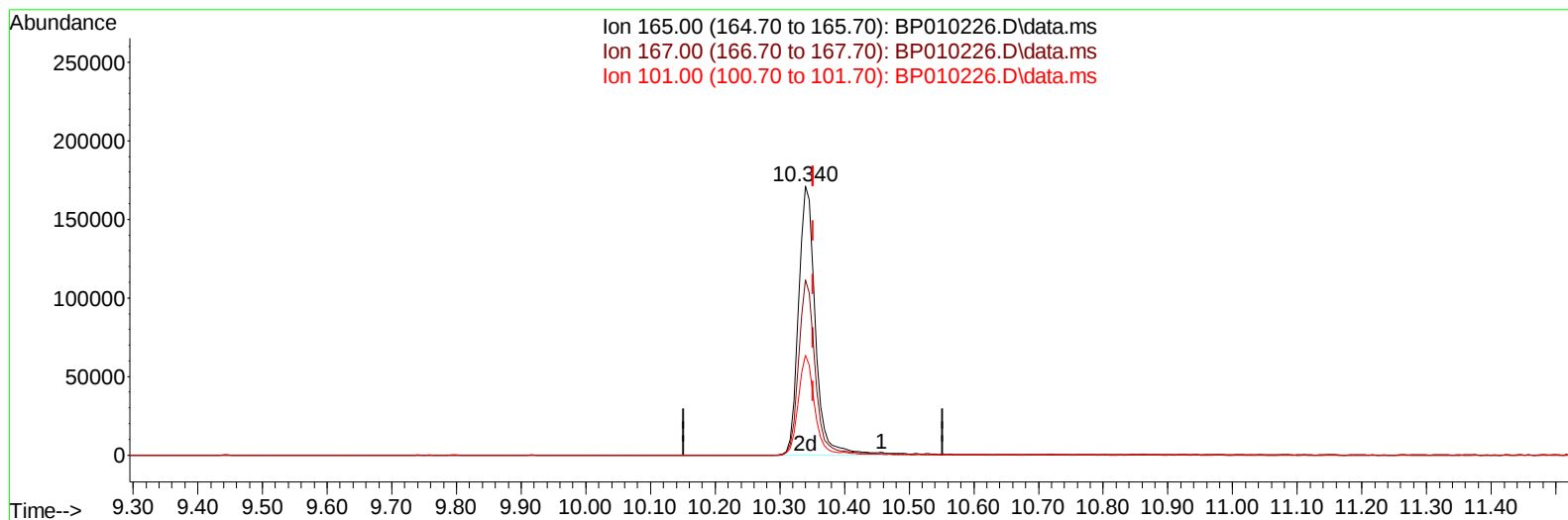
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TIC: BP010226.D\data.ms

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

10.340min (-0.012) 27.38 ng/ul m

response 313951

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	65.21
101.00	49.80	37.15#
0.00	0.00	0.00

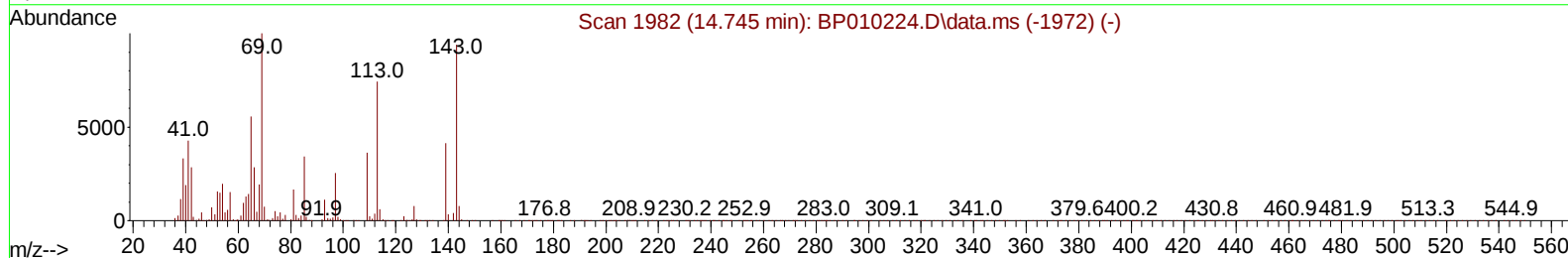
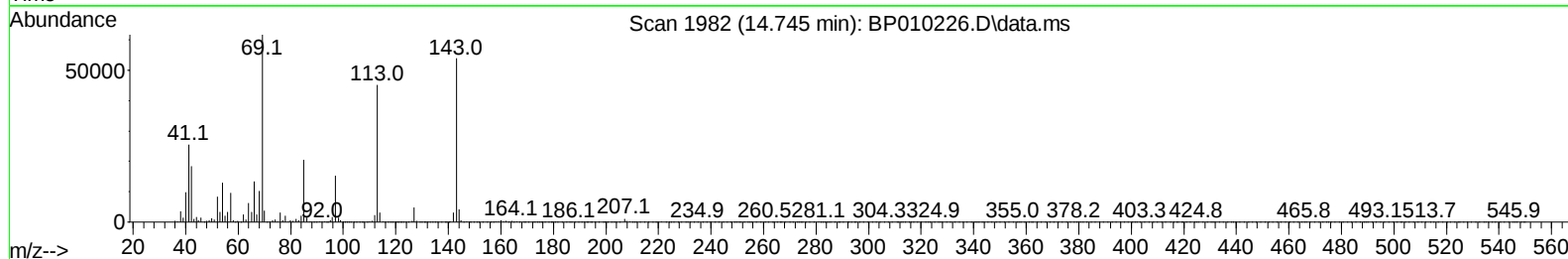
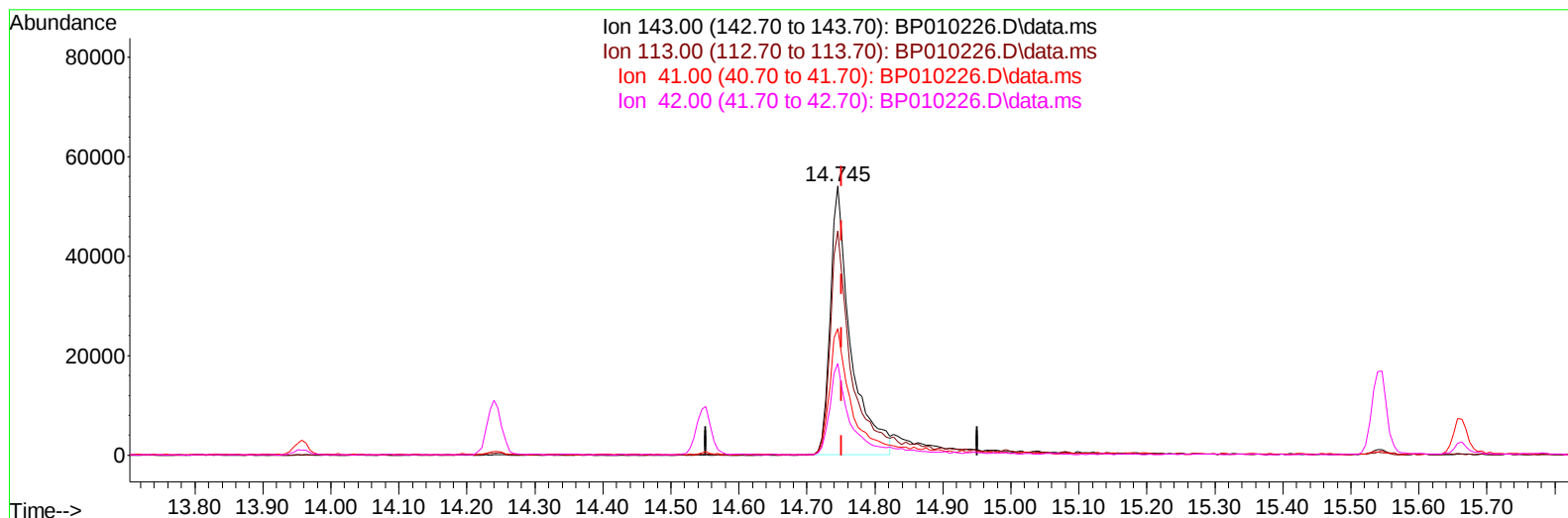
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TIC: BP010226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 17.61 ng/ul

response 113957

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.44#
41.00	23.20	47.04#
42.00	18.00	34.00#

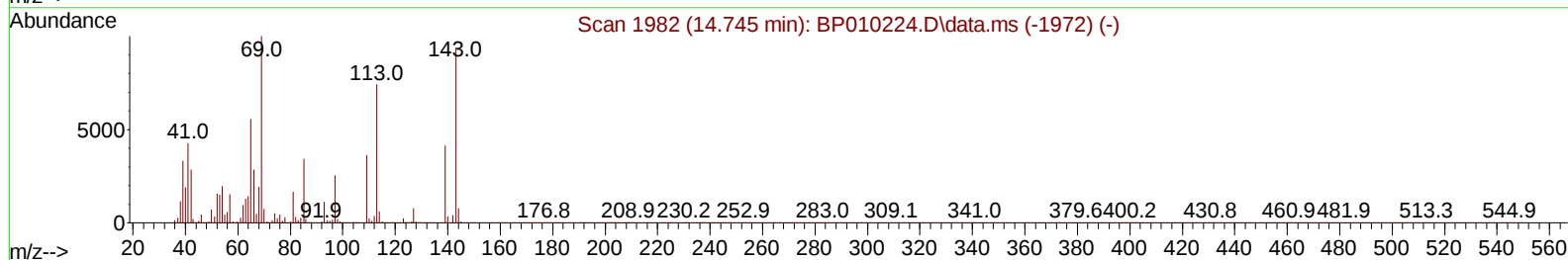
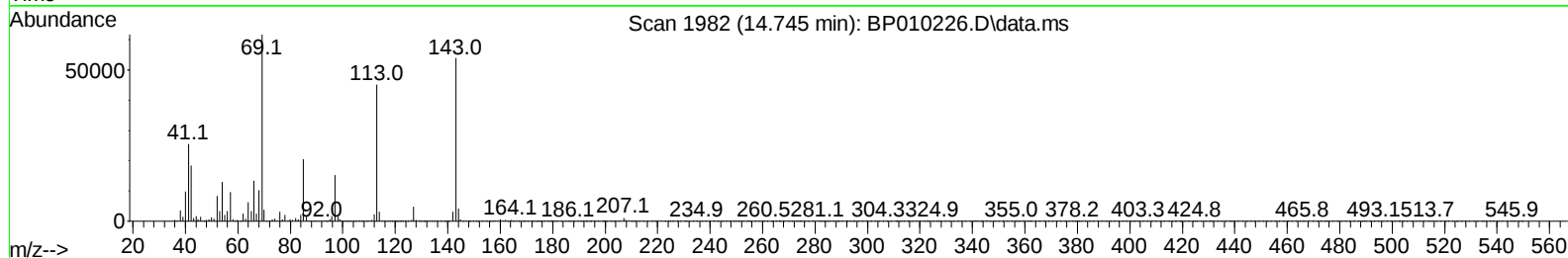
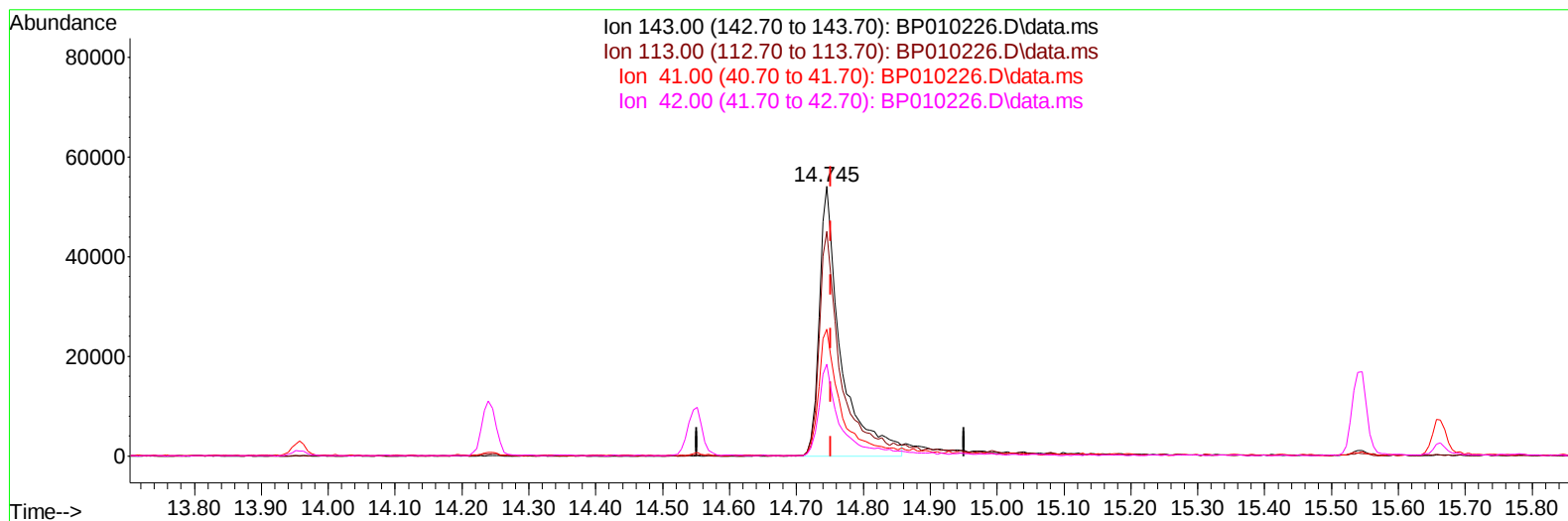
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
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TIC: BP010226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 18.69 ng/ul m

response 120974

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.44#
41.00	23.20	47.04#
42.00	18.00	34.00#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	153491	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	678163	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	451993	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	872675	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	603514	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	560918	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	21639	6.148	ng/uL	0.00
4) Pyridine-d5	3.775	84	246058	24.849	ng/ul	0.00
7) Phenol-d5	7.081	99	375462	26.893	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	268815	32.088	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	312319	30.206	ng/ul	-0.01
15) 4-Methylphenol-d8	8.622	113	323651	27.980	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	165582	30.699	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	176622	29.983	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	313951m	27.377	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	313967	19.094	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1096602	31.189	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1239683	29.133	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.745	143	120974m	18.689	ng/ul	0.00
60) Fluorene-d10	15.545	176	919571	30.404	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	113929	20.647	ng/ul	0.00
73) Anthracene-d10	17.404	188	1348894	31.950	ng/ul	0.00
81) Pyrene-d10	19.633	212	1345179	39.193	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	989581	33.652	ng/ul	-0.01

Target Compounds Qvalue

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

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