Quantitation Report (Qedit)

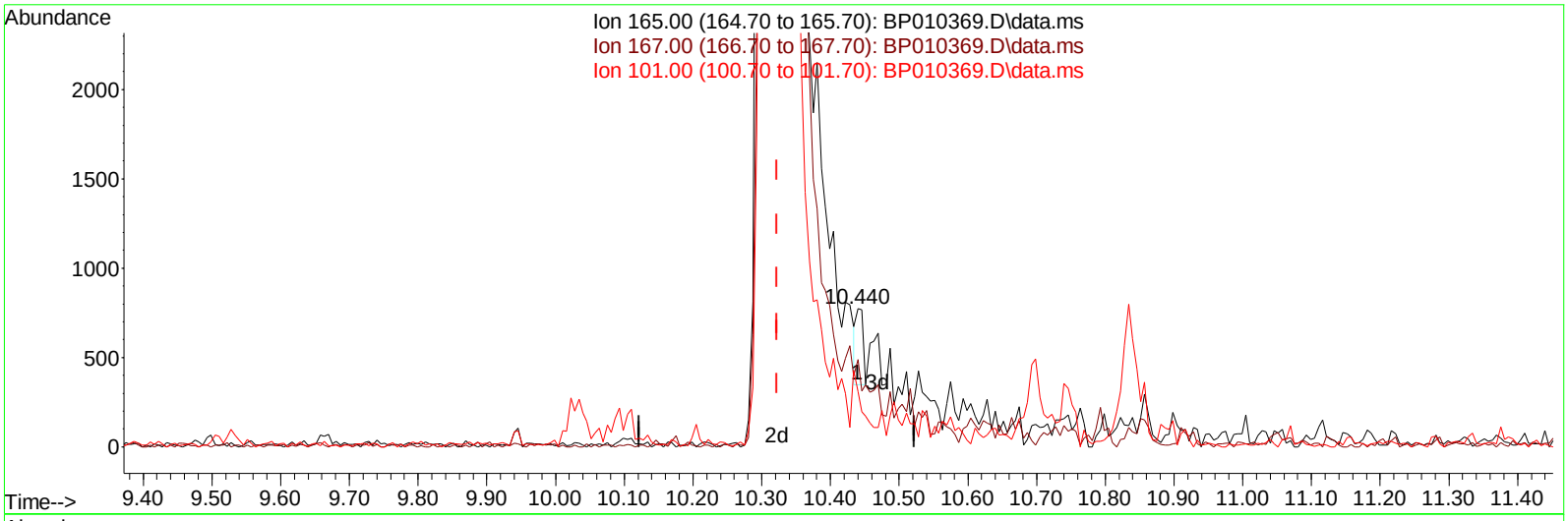
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010369.D
Acq On : 16 May 2022 22:02
Operator : CG/JU
Sample : N2801-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

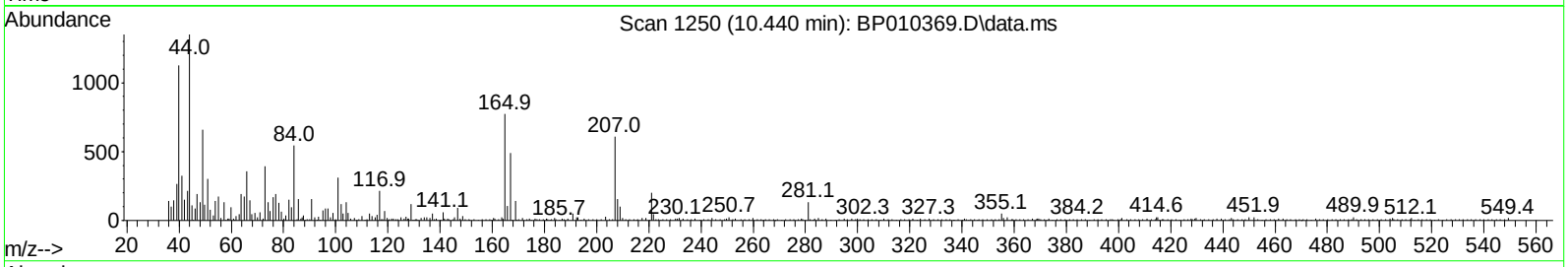
Quant Time: May 16 23:22:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

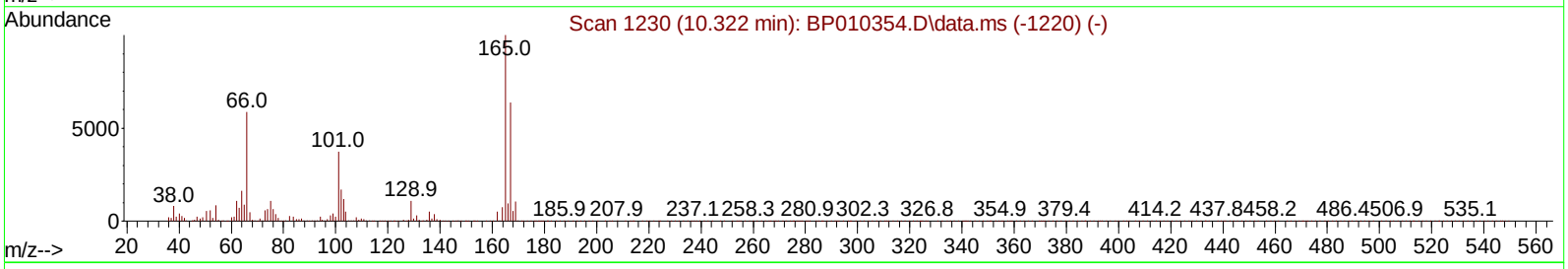
Instrument :
BNA_P
ClientSampleId :
EXLB9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/17/2022
Supervised By :mohammad ahmed 05/18/2022







TIC: BP010369.D\data.ms

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

10.440min (+ 0.118) 0.03 ng/ul

response 297

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.39
101.00	49.80	40.10
0.00	0.00	0.00

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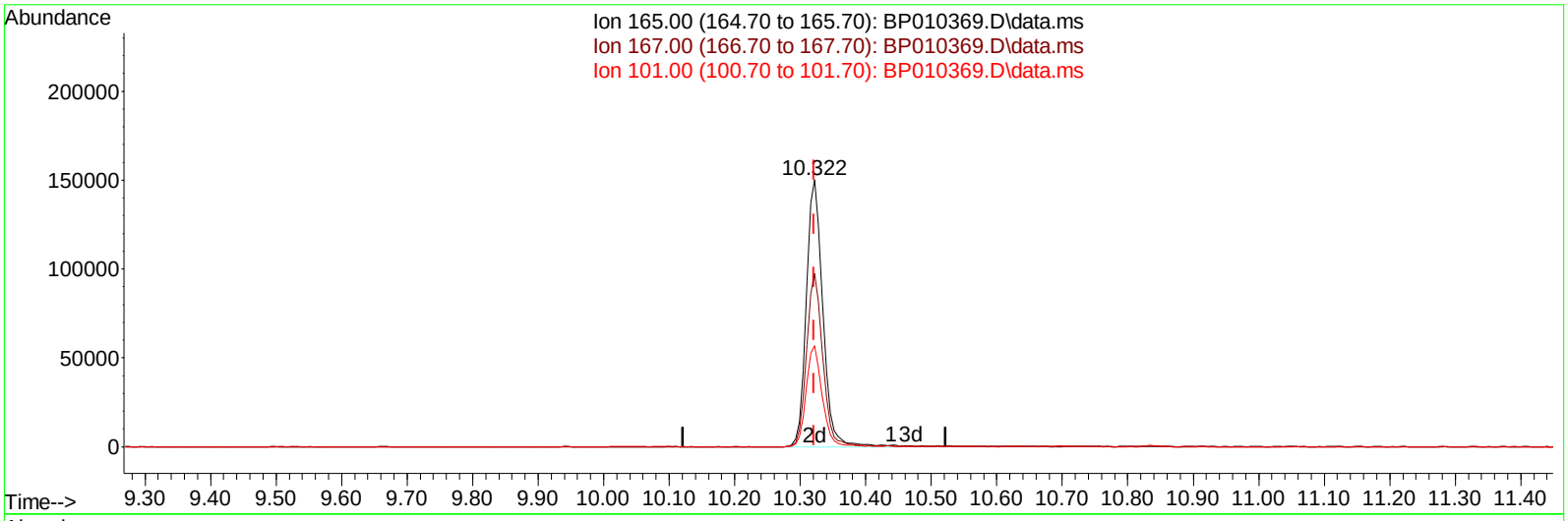

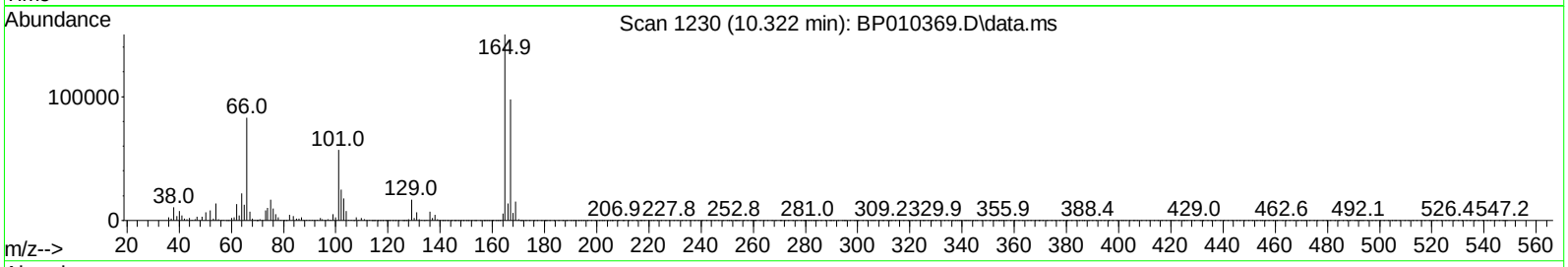

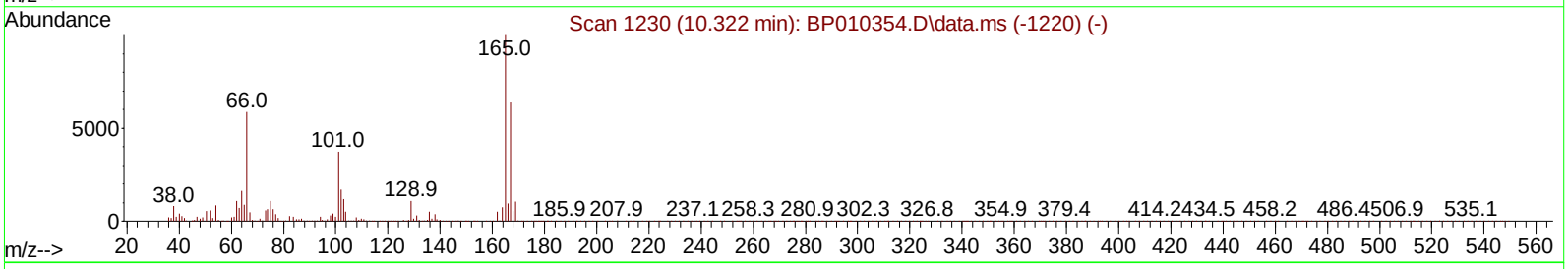

TIC: BP010369.D\data.ms

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

10.322min (+ 0.000) 30.86 ng/ul m

response 264576

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	65.05
101.00	49.80	37.84#
0.00	0.00	0.00

Quantitation Report (Qedit)

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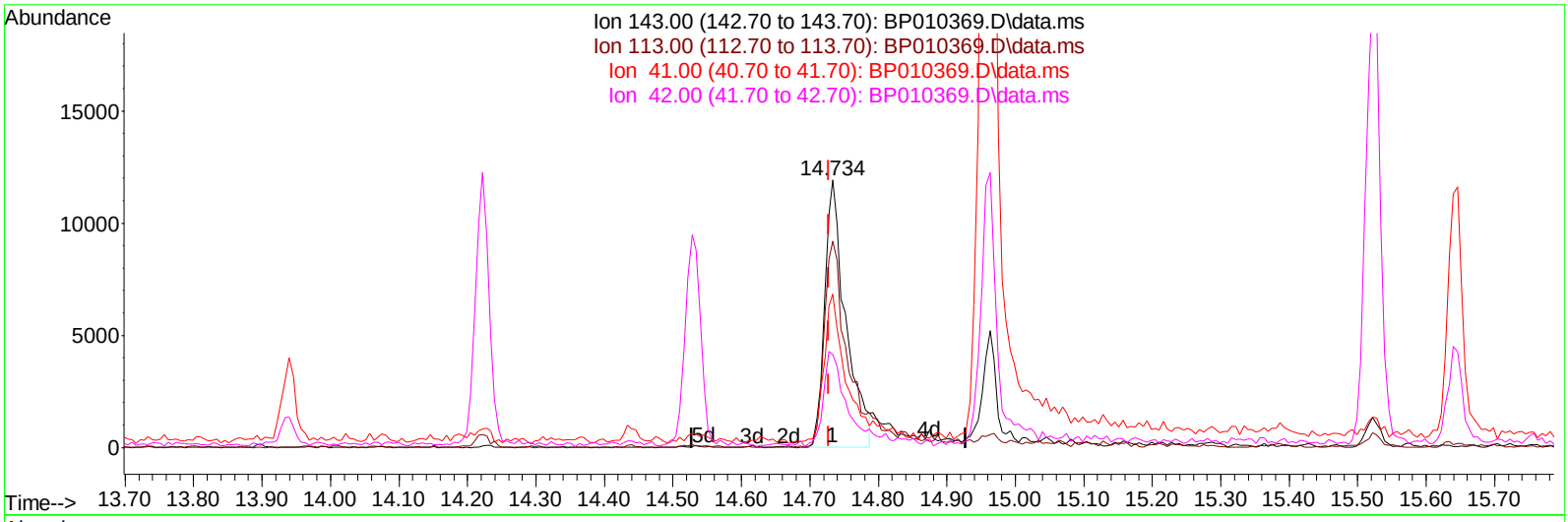

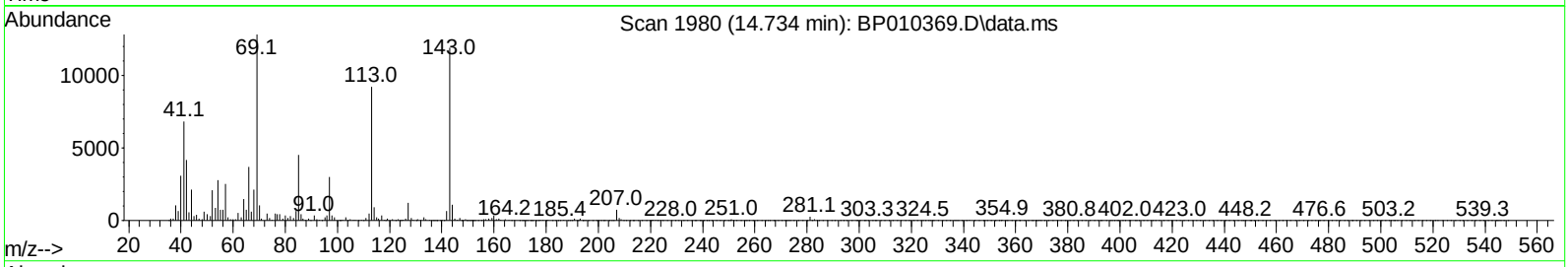

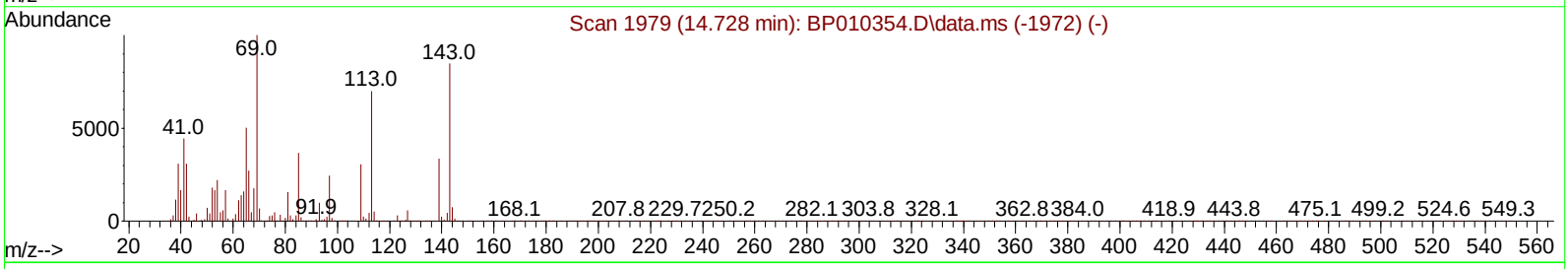

TIC: BP010369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 5.01 ng/ul

response 25424

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.30#
41.00	23.20	57.35#
42.00	18.00	34.87#

Quantitation Report (Qedit)

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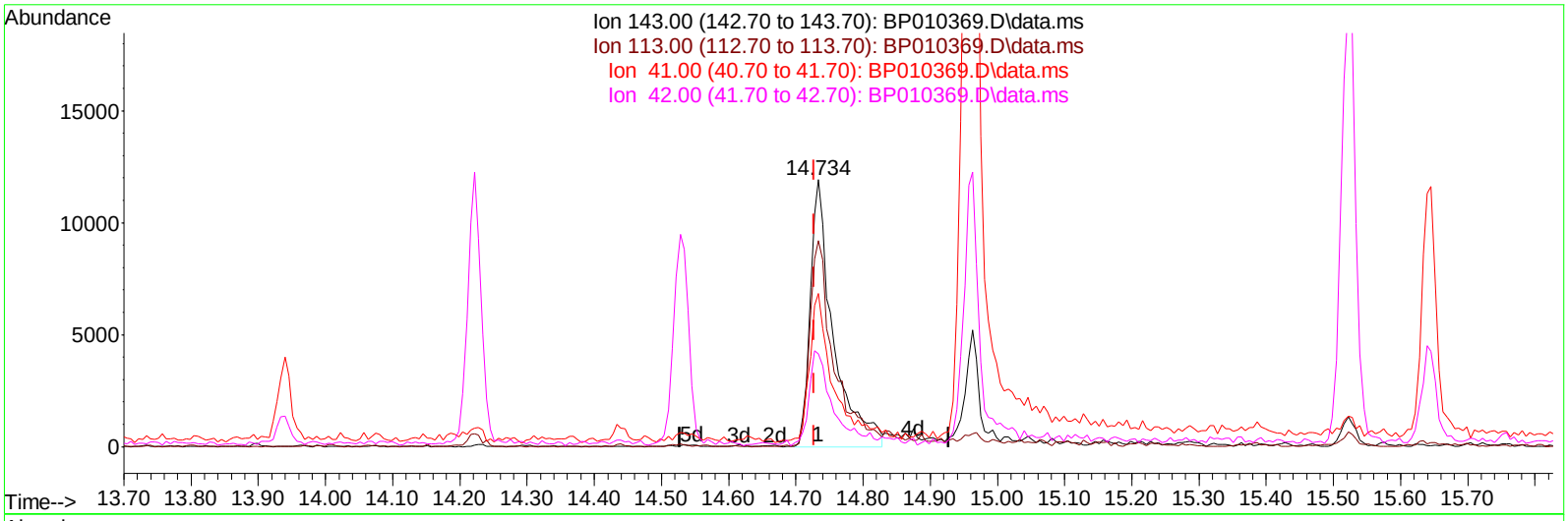

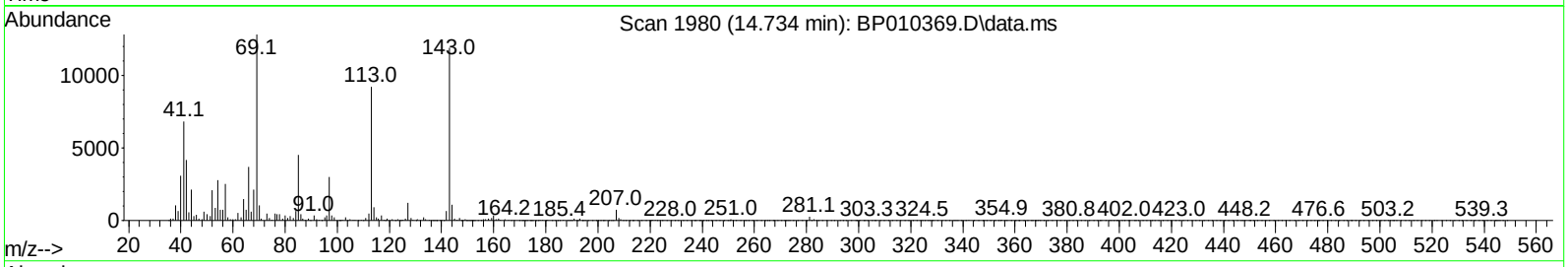

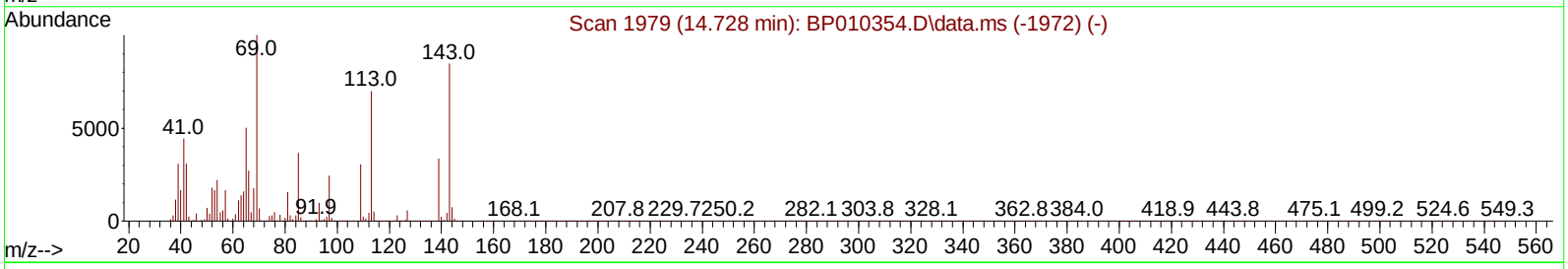

TIC: BP010369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 5.55 ng/ul m

response 28169

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.30#
41.00	23.20	57.35#
42.00	18.00	34.87#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/17/2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	126624	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	552792	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	387271	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	863281	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	922056	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	890729	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	6803	2.290	ng/uL	0.00
4) Pyridine-d5	3.770	84	85888	10.182	ng/ul	0.00
7) Phenol-d5	7.064	99	67532	6.354	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	246621	34.538	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	209610	25.775	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	144123	15.957	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	152688	35.689	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	153541	33.865	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	264576m	30.862	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	379097	29.740	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1120878	39.284	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1199182	36.436	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	28169m	5.547	ng/ul	0.00
60) Fluorene-d10	15.522	176	973549	39.219	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	162833	32.252	ng/ul	0.00
73) Anthracene-d10	17.381	188	1714727	43.865	ng/ul	0.00
81) Pyrene-d10	19.616	212	2157511	44.060	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2019387	45.149	ng/ul	0.00
Target Compounds						Qvalue

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

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