Quantitation Report (Qedit)

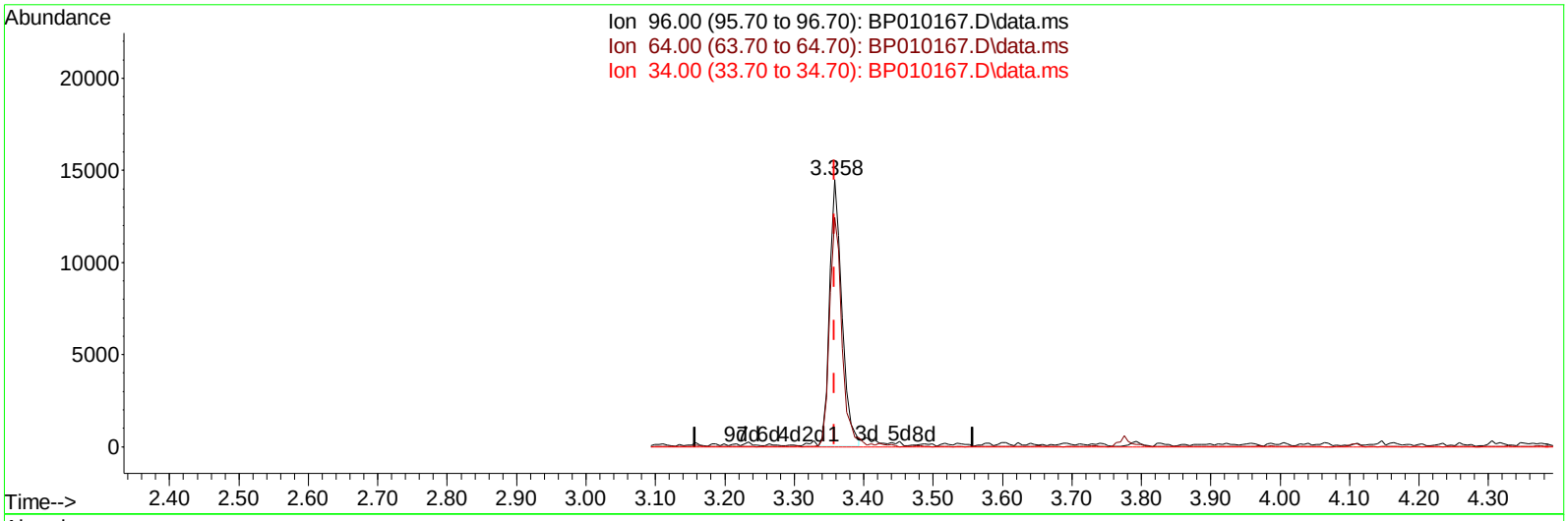
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010167.D
Acq On : 02 May 2022 11:48
Operator : CG/JU
Sample : PB144495BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

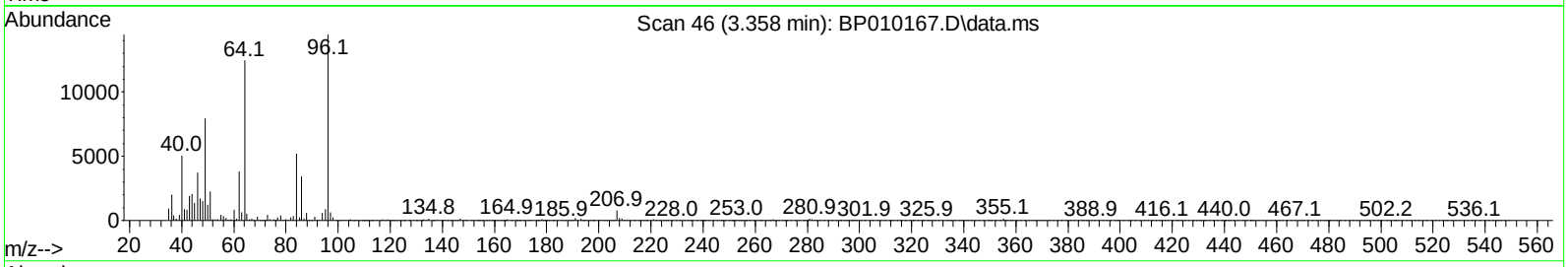
Quant Time: May 03 01:06:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

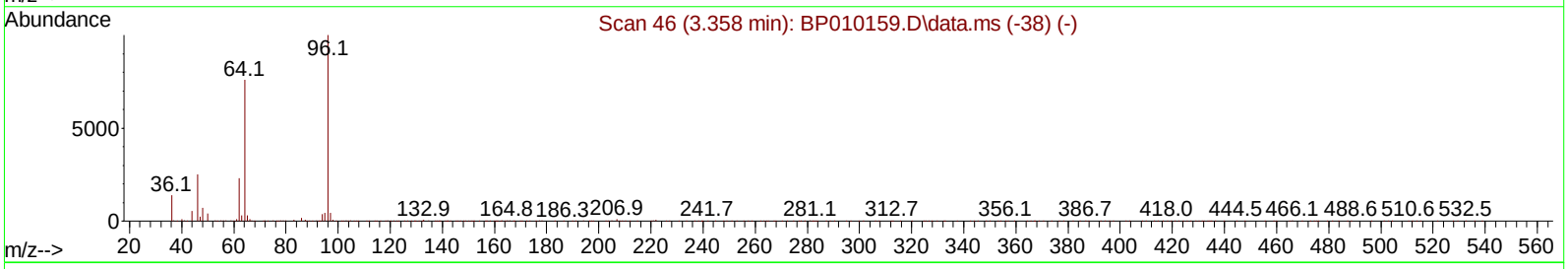
Instrument :
BNA_P
ClientSampleId :
SBLK495

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/03/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010167.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.358min (+ 0.000) 5.75 ng/uL

response 18076

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	85.95#
34.00	0.00	0.00
0.00	0.00	0.00

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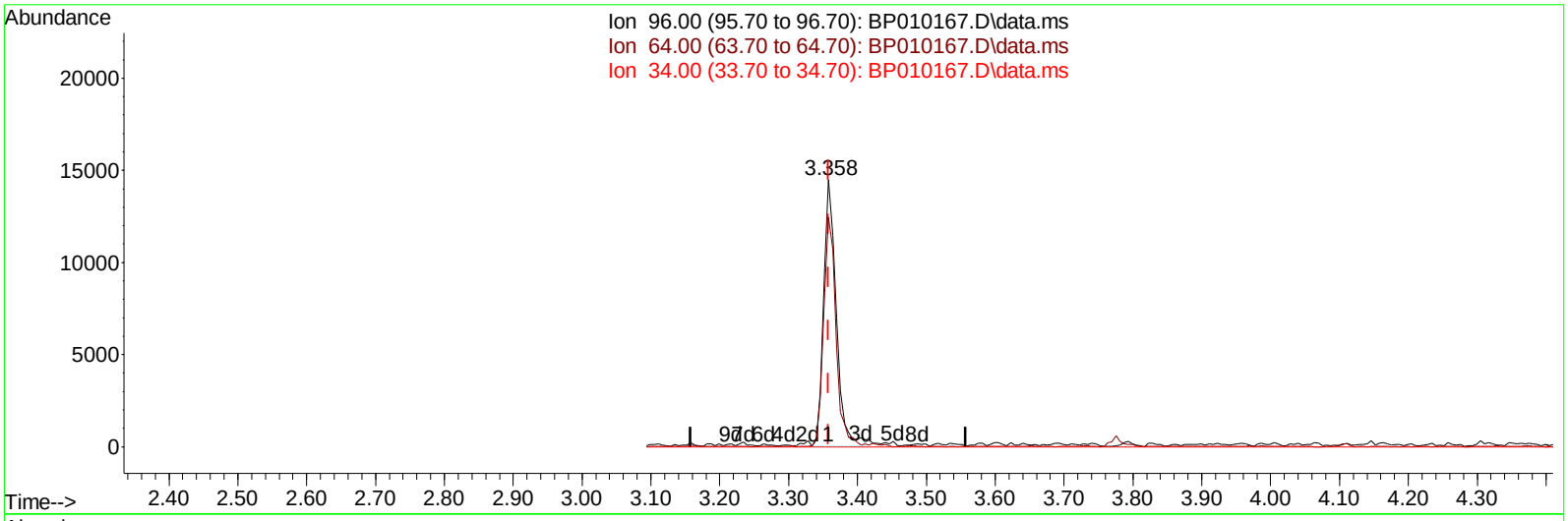

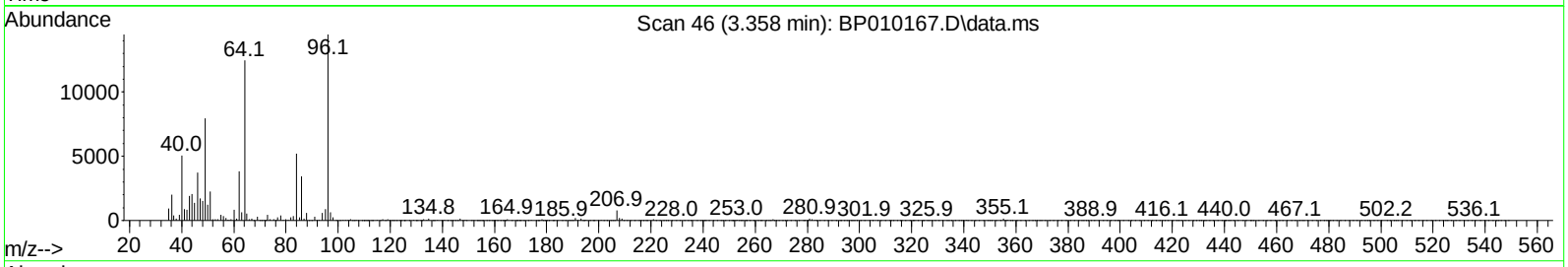

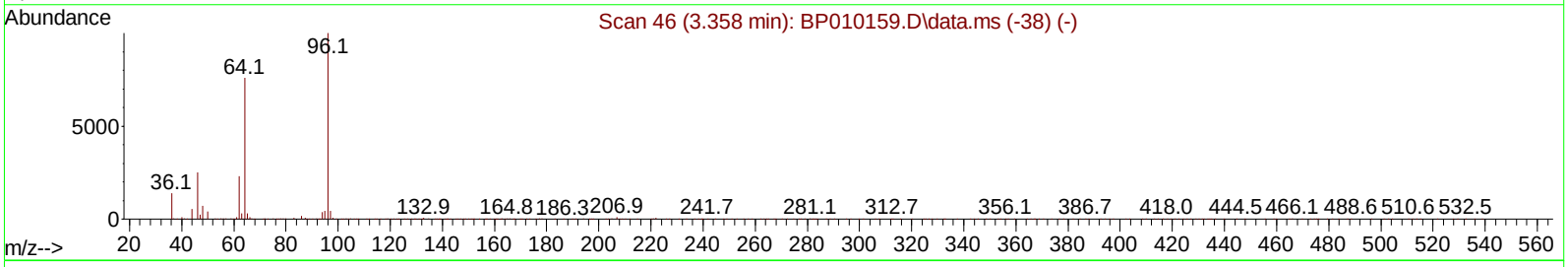

TIC: BP010167.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.358min (+ 0.000) 5.89 ng/uL m

response 18514

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	85.95#
34.00	0.00	0.00
0.00	0.00	0.00

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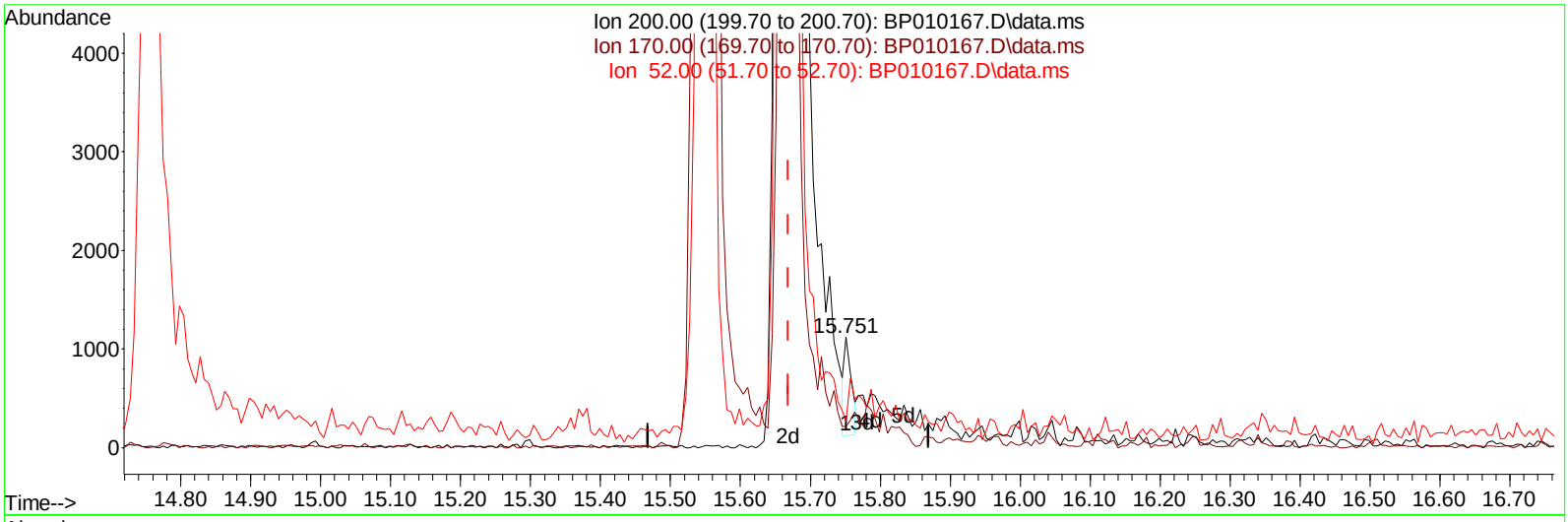

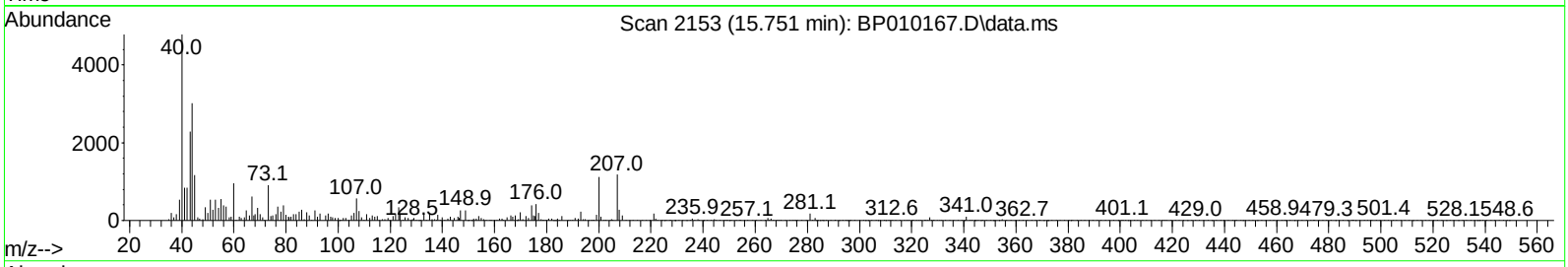

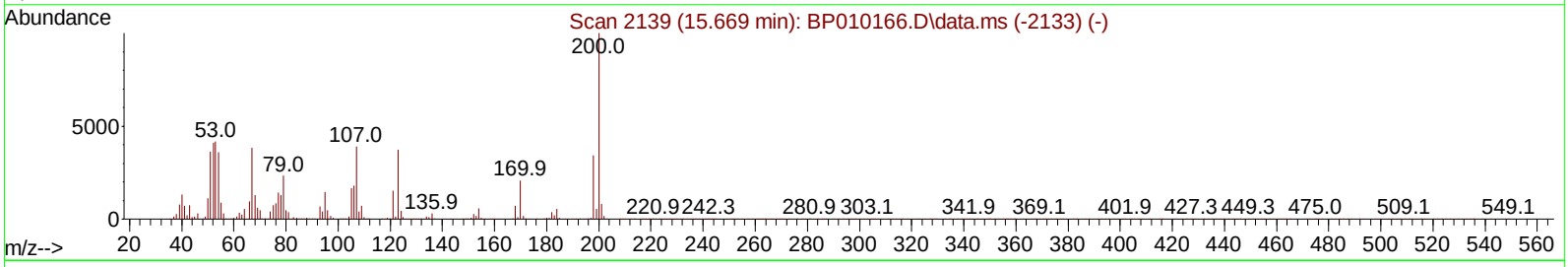

TIC: BP010167.D\data.ms

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

15.751min (+ 0.082) 0.14 ng/ul

response 689

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	19.07
52.00	21.20	23.98
0.00	0.00	0.00

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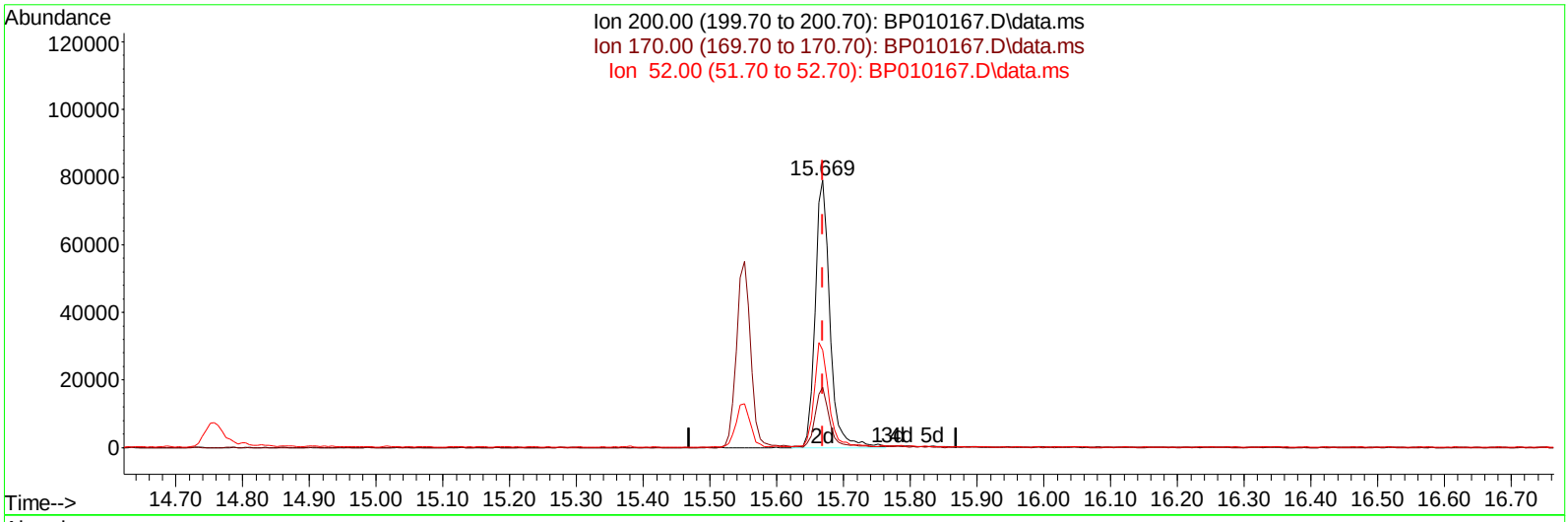

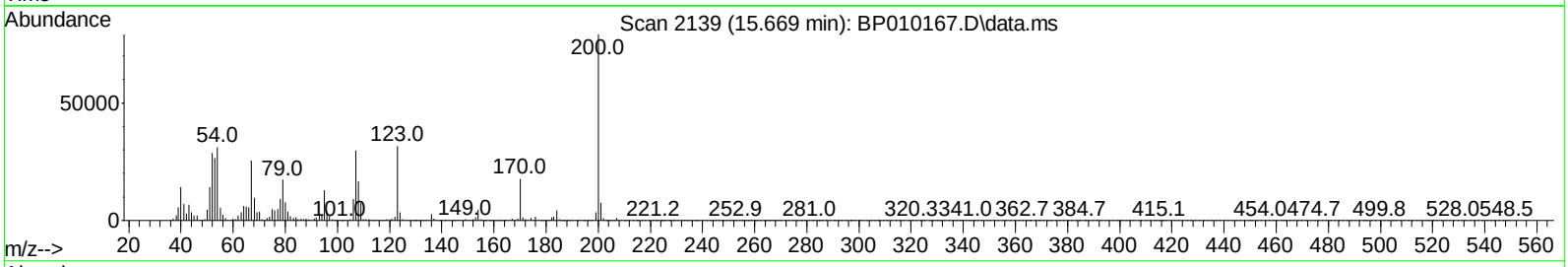

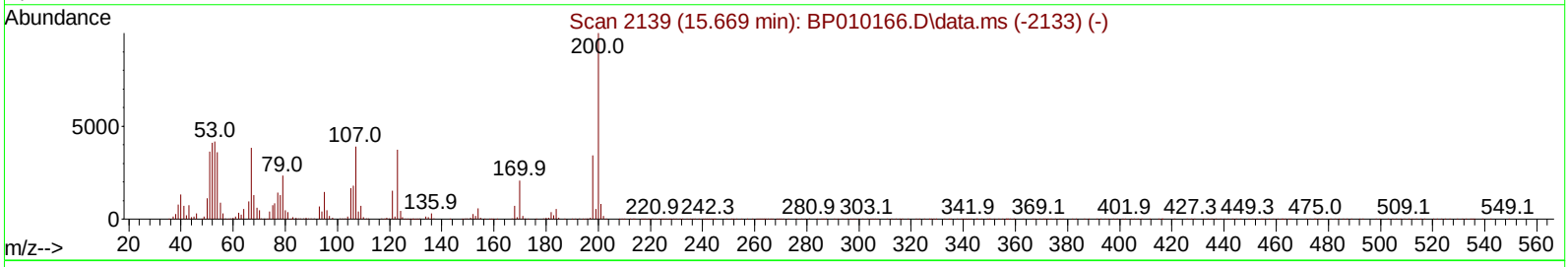

TIC: BP010167.D\data.ms

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

15.669min (+ 0.000) 25.31 ng/ul m

response 120786

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	22.51
52.00	21.20	36.33#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 05/05/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.922	152	137083	20.000	ng/ul		0.00
20) Naphthalene-d8	10.728	136	560374	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.557	164	362374	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.310	188	754602	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.392	240	705172	20.000	ng/ul		0.00
88) Perylene-d12	23.839	264	674784	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.358	96	18514m	5.889	ng/uL		0.00
4) Pyridine-d5	3.775	84	245040	27.708	ng/ul		0.00
7) Phenol-d5	7.087	99	308444	24.737	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	217218	29.032	ng/ul		0.00
11) 2-Chlorophenol-d4	7.458	132	258304	27.972	ng/ul		0.00
15) 4-Methylphenol-d8	8.634	113	260996	25.264	ng/ul		0.00
21) Nitrobenzene-d5	9.081	128	132025	29.623	ng/ul		0.00
24) 2-Nitrophenol-d4	9.805	143	147186	30.238	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.352	165	245729	25.932	ng/ul		0.00
31) 4-Chloroaniline-d4	10.863	131	345749	25.447	ng/ul		0.00
46) Dimethylphthalate-d6	13.963	166	844281	29.951	ng/ul		0.00
49) Acenaphthylene-d8	14.251	160	938587	27.512	ng/ul		0.00
54) 4-Nitrophenol-d4	14.757	143	119686	23.063	ng/ul		0.00
60) Fluorene-d10	15.551	176	708543	29.221	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	120786m	25.314	ng/ul		0.00
73) Anthracene-d10	17.404	188	1101816	30.181	ng/ul		0.00
81) Pyrene-d10	19.639	212	1299792	32.411	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.680	264	1095347	30.963	ng/ul		0.00

Target Compounds Qvalue

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

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