Quantitation Report (Qedit)

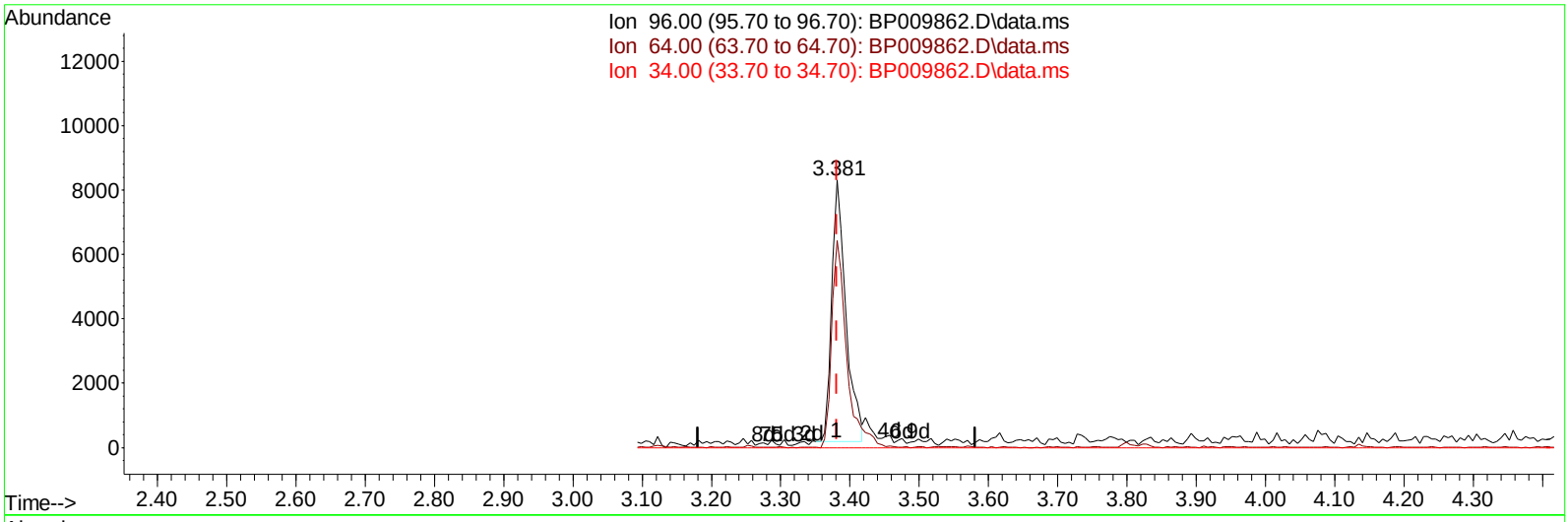
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
Data File : BP009862.D
Acq On : 15 Apr 2022 15:37
Operator : CG/JU
Sample : N2422-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

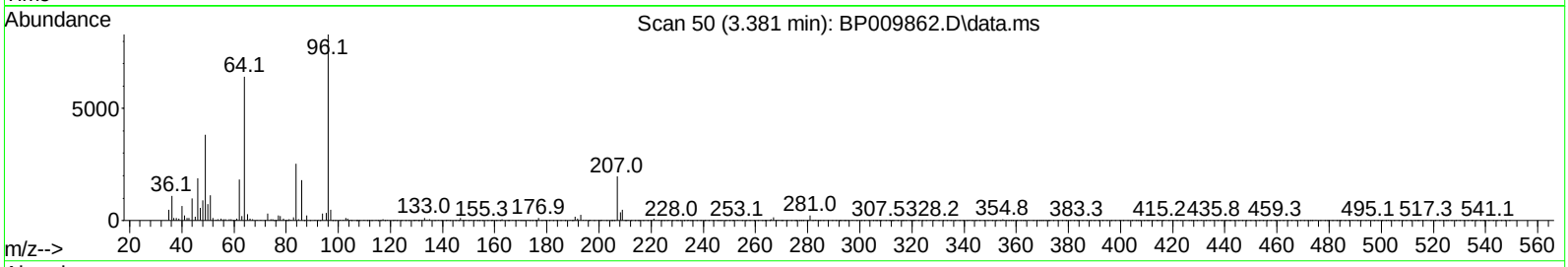
Quant Time: Apr 18 01:01:45 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 18 00:59:30 2022
Response via : Initial Calibration

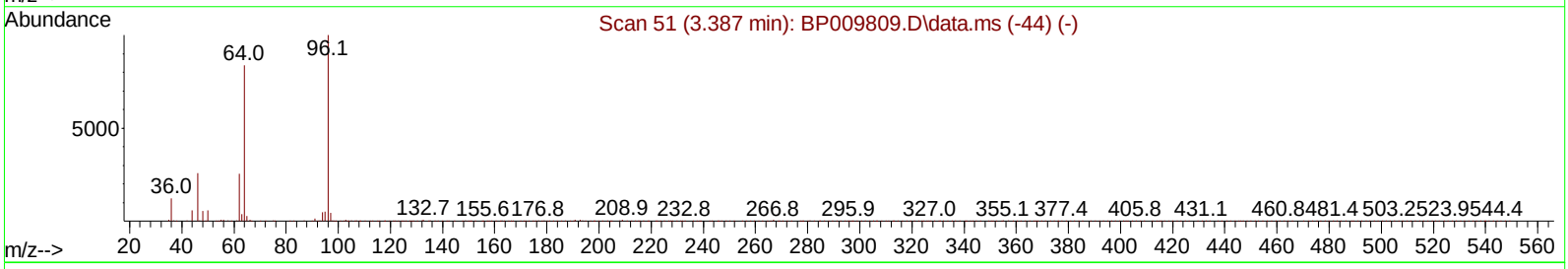
Instrument :
BNA_P
ClientSampleId :
C0J57

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022
Supervised By :Yogesh Patel 04/18/2022







TIC: BP009862.D\data.ms

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

3.381min (+ 0.000) 4.56 ng/uL

response 11506

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	77.37#
34.00	0.00	0.00
0.00	0.00	0.00

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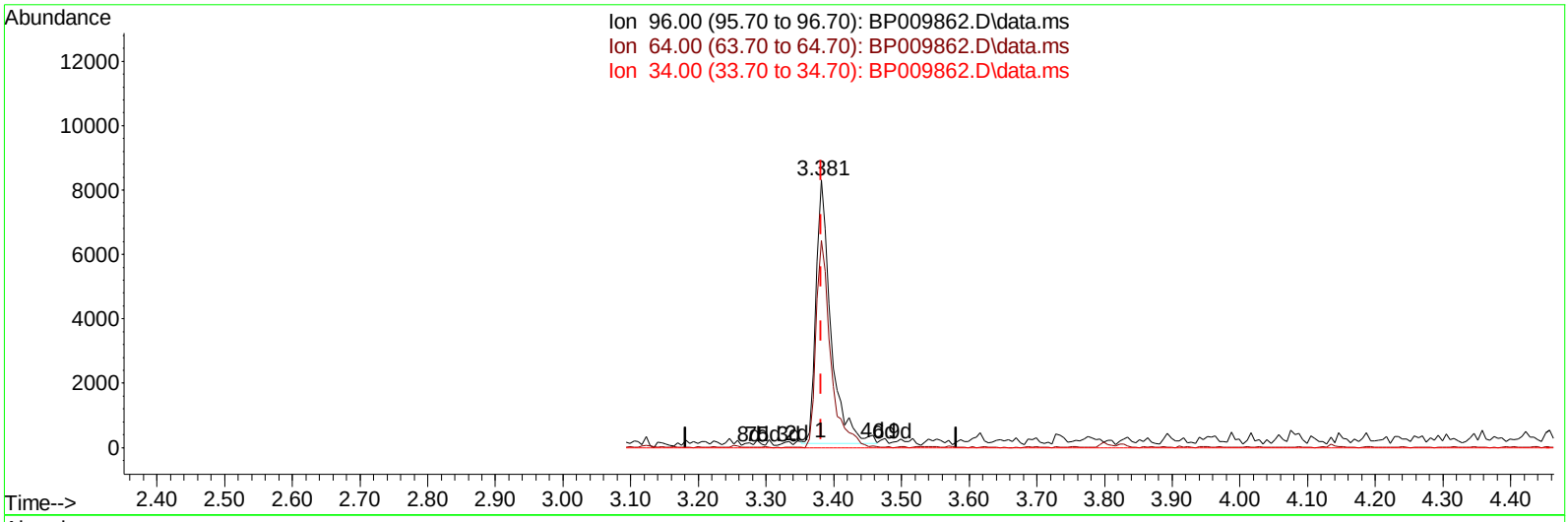

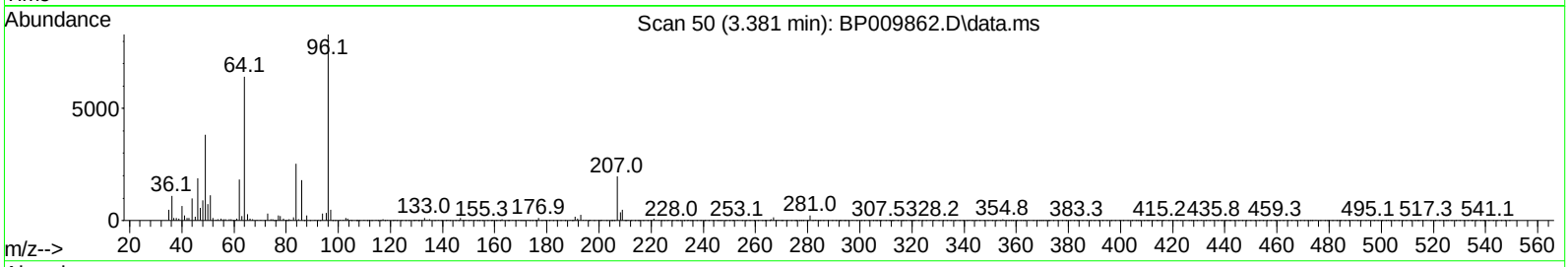

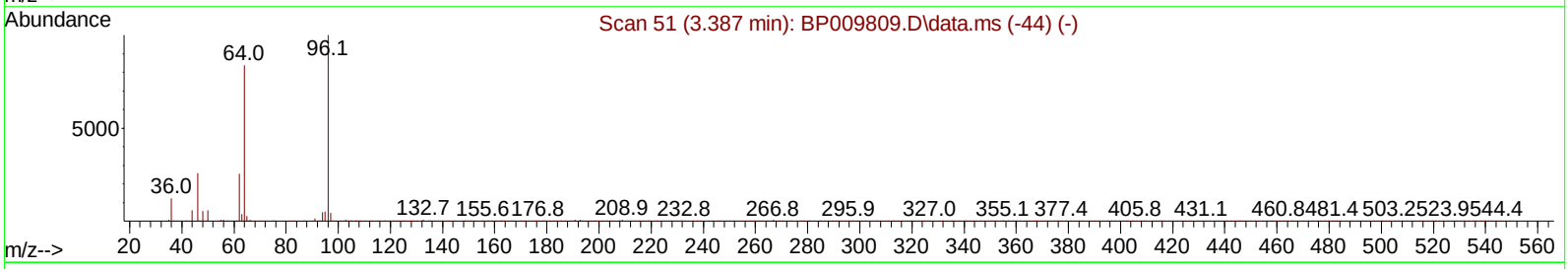

TIC: BP009862.D\data.ms

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

3.381min (+ 0.000) 4.97 ng/uL m

response 12553

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	77.37#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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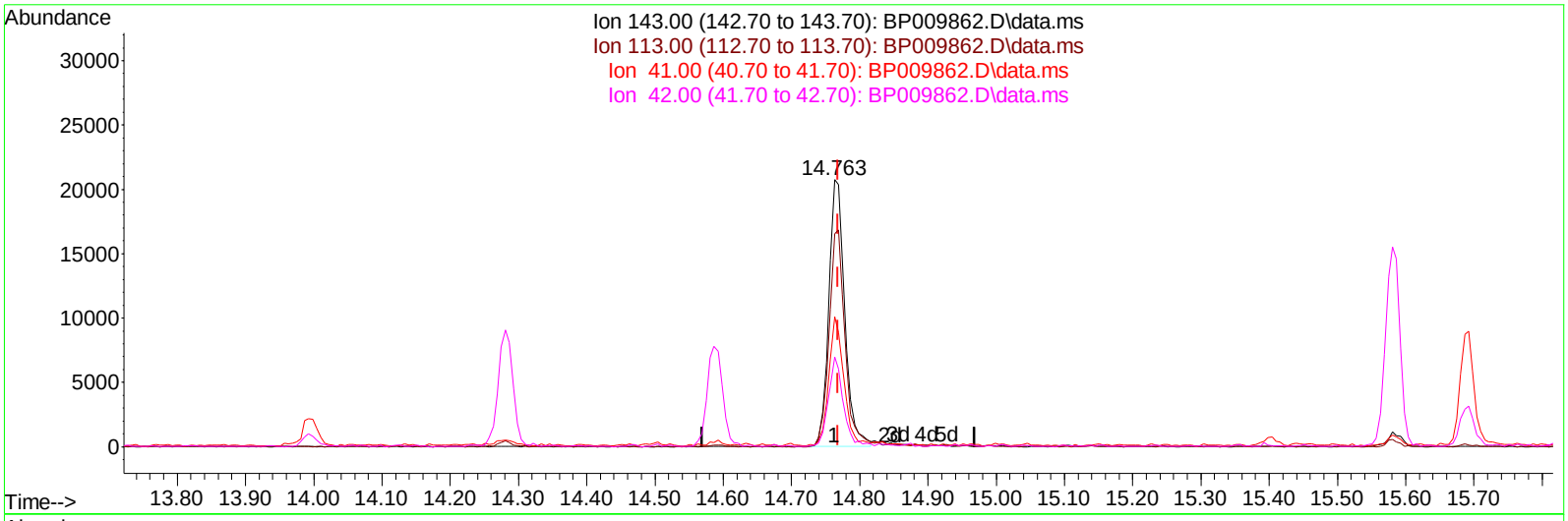

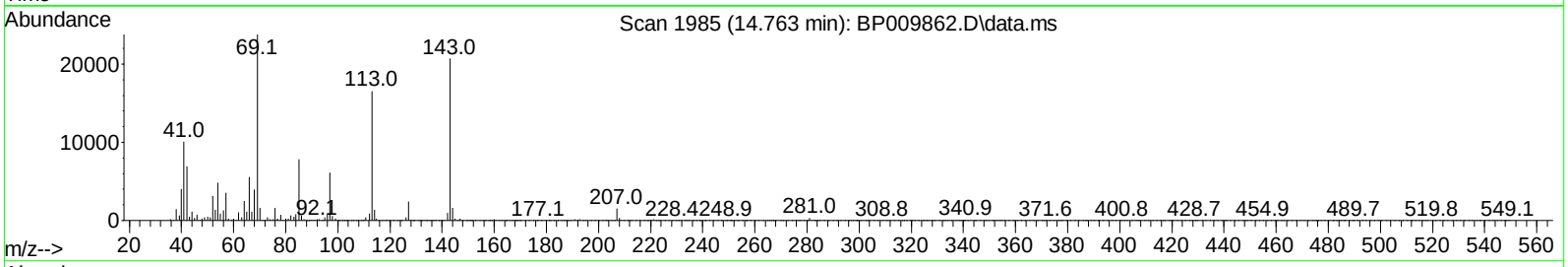

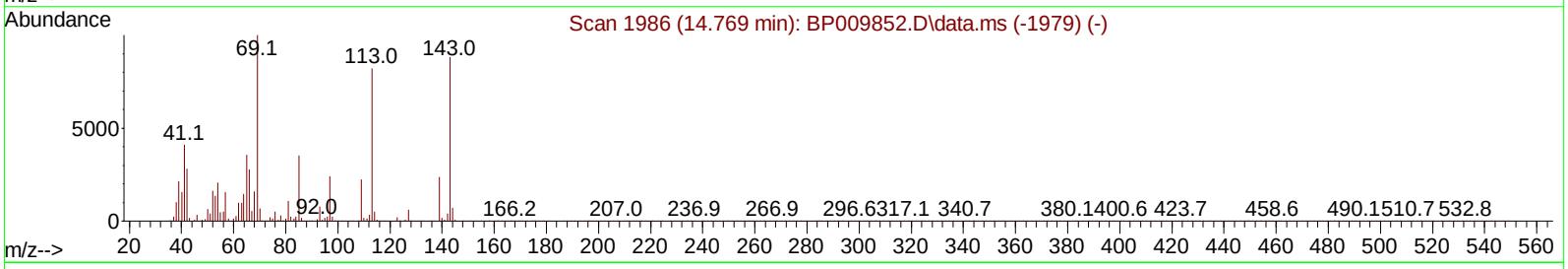

TIC: BP009862.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 6.70 ng/ul

response 33723

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.87#
41.00	23.20	48.62#
42.00	18.00	33.54#

Quantitation Report (Qedit)

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ClientSampleId :
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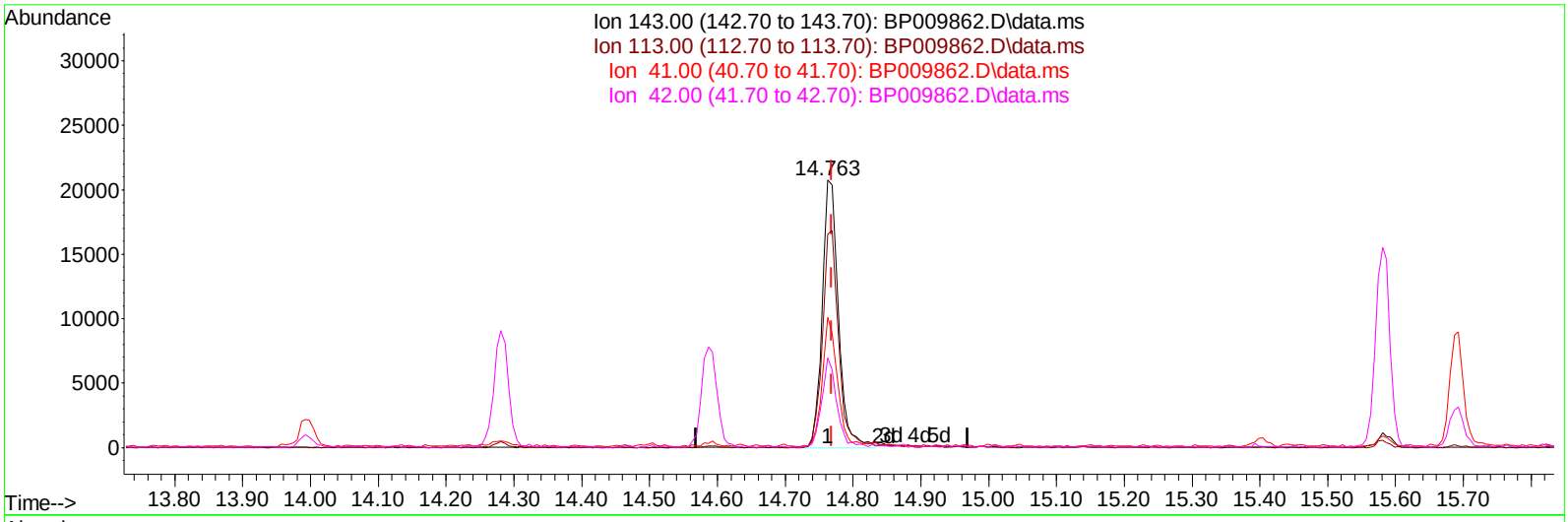

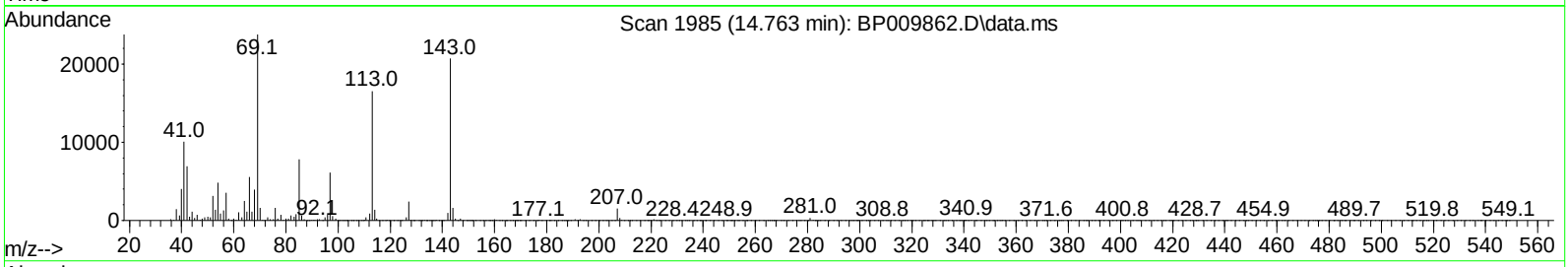

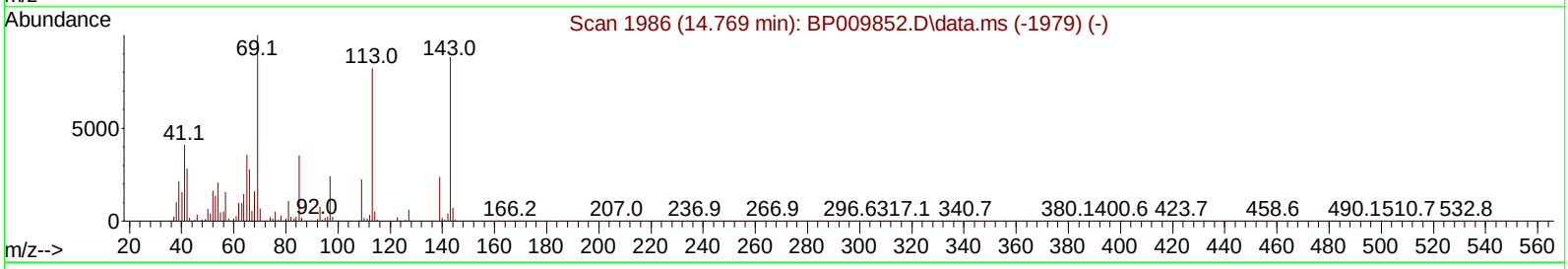

TIC: BP009862.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 6.77 ng/ul m

response 34100

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.87#
41.00	23.20	48.62#
42.00	18.00	33.54#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 04/18/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	112719	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	497634	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	353350	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	804644	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	861267	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	832417	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12553m	4.970	ng/uL	0.00
4) Pyridine-d5	3.805	84	54409	7.813	ng/ul	0.00
7) Phenol-d5	7.110	99	68614	6.933	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	187296	31.791	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	180825	24.224	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	132666	16.059	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	120145	33.480	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	122096	31.033	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	212691	25.834	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	316754	27.013	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	928823	33.464	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	996362	30.300	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	34100m	6.774	ng/ul	0.00
60) Fluorene-d10	15.581	176	816378	34.963	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	137888	28.927	ng/ul	0.00
73) Anthracene-d10	17.439	188	1430815	37.139	ng/ul	0.00
81) Pyrene-d10	19.669	212	1759437	37.404	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1697499	39.159	ng/ul	0.00
Target Compounds						Qvalue

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

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