Quantitation Report (Qedit)

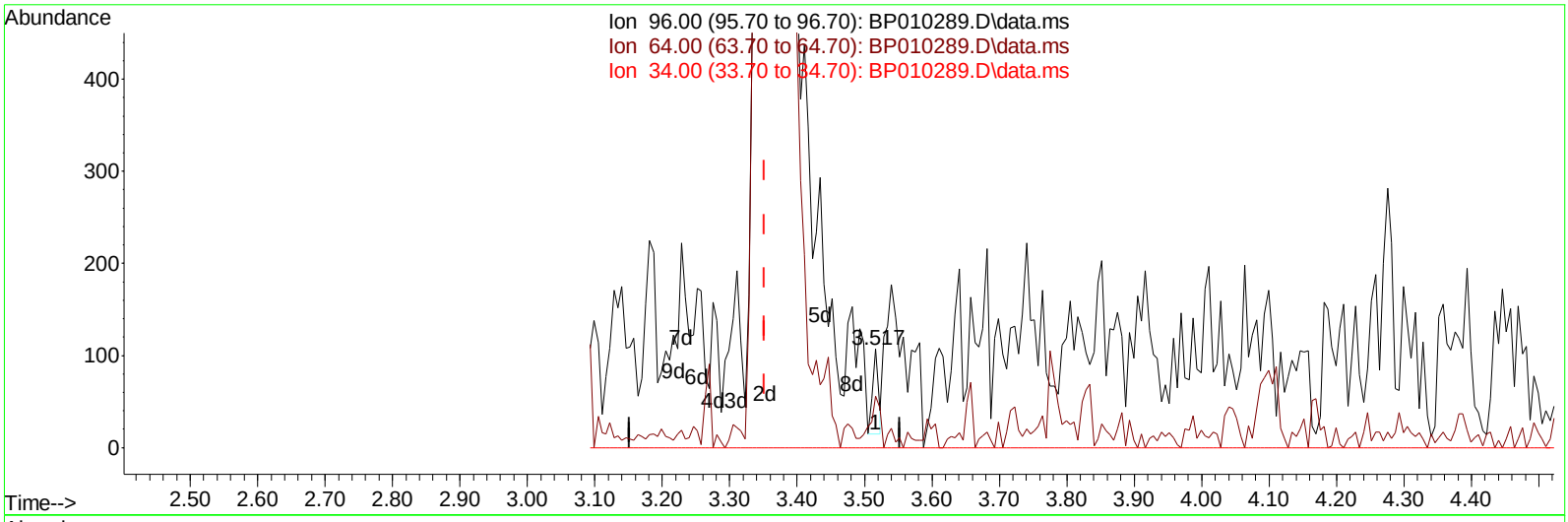
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010289.D
Acq On : 11 May 2022 15:24
Operator : CG/JU
Sample : N2632-13
Misc :
ALS Vial : 11 Sample Multiplier: 1

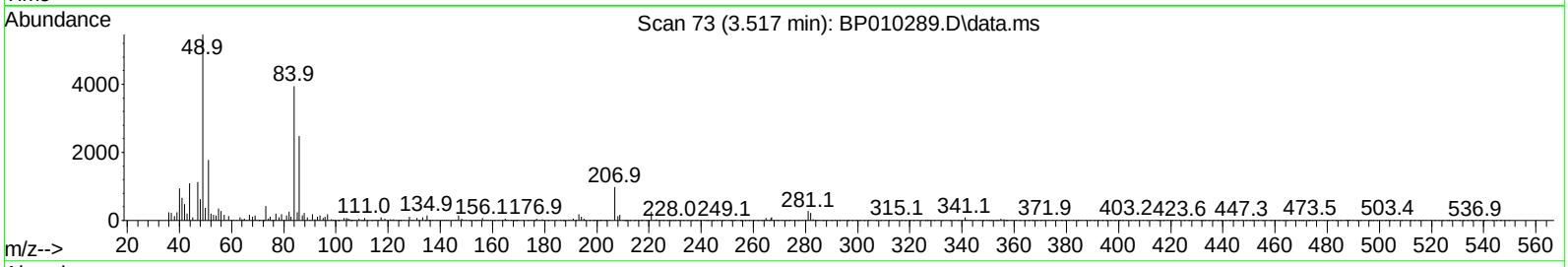
Quant Time: May 12 01:13:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

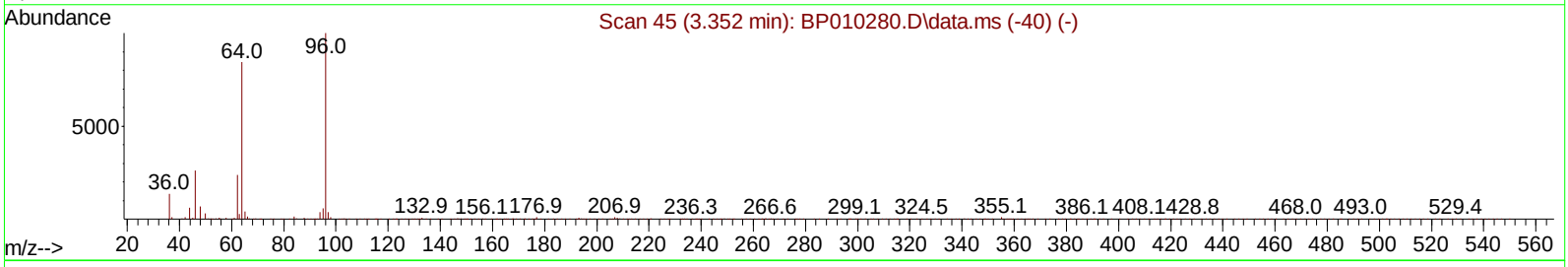
Instrument :
BNA_P
ClientSampleId :
EW4B5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022







TIC: BP010289.D\data.ms

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

3.517min (+ 0.165) 0.02 ng/uL

response 55

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	52.34
34.00	0.00	0.00
0.00	0.00	0.00

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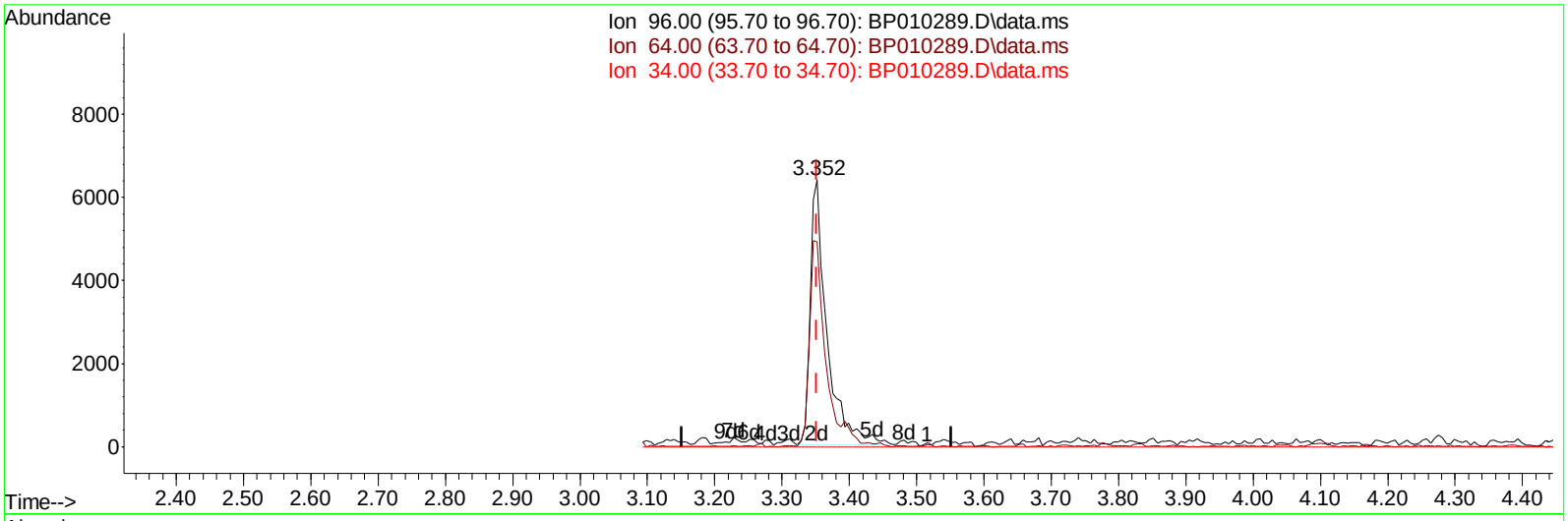

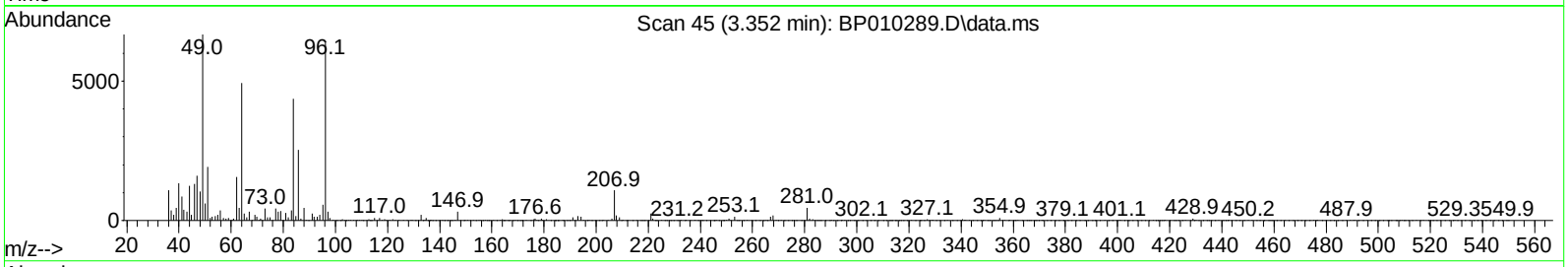

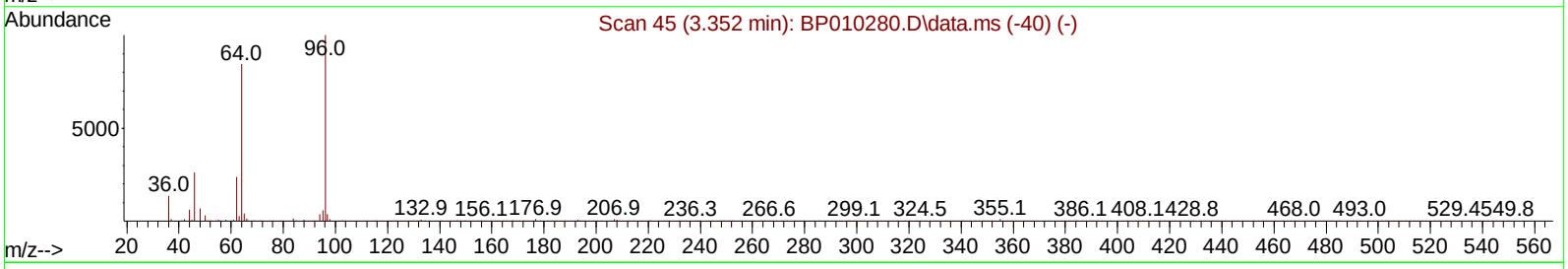

TIC: BP010289.D\data.ms

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

3.352min (+ 0.000) 3.59 ng/uL m

response 11101

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	76.76#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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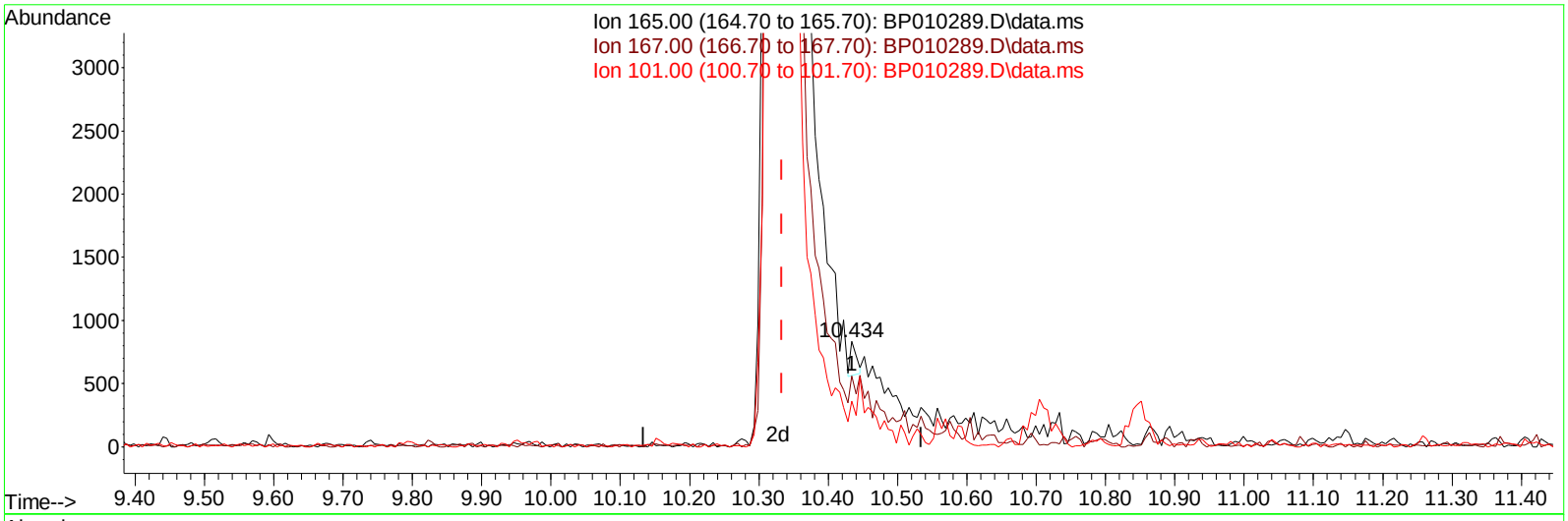

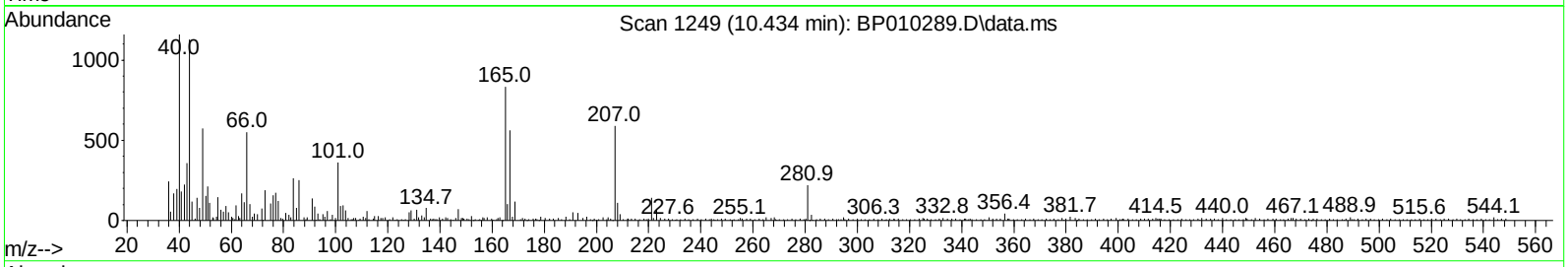

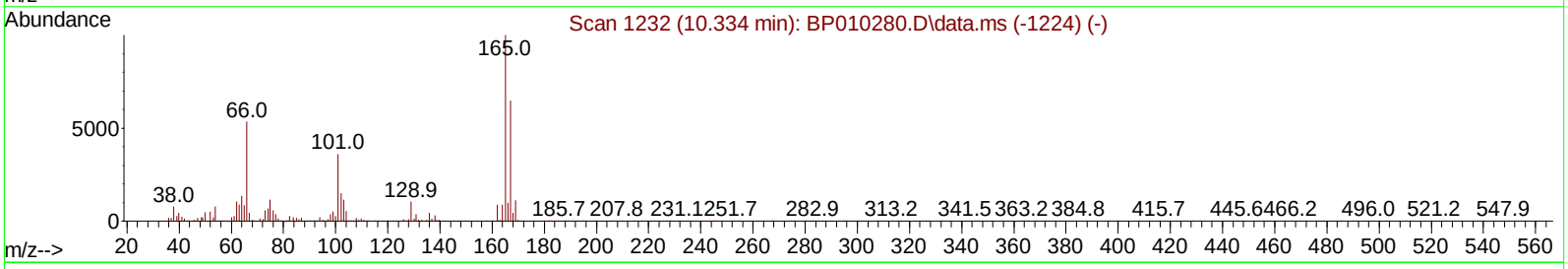

TIC: BP010289.D\data.ms

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

10.434min (+ 0.100) 0.02 ng/ul

response 168

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	67.23
101.00	49.80	43.22
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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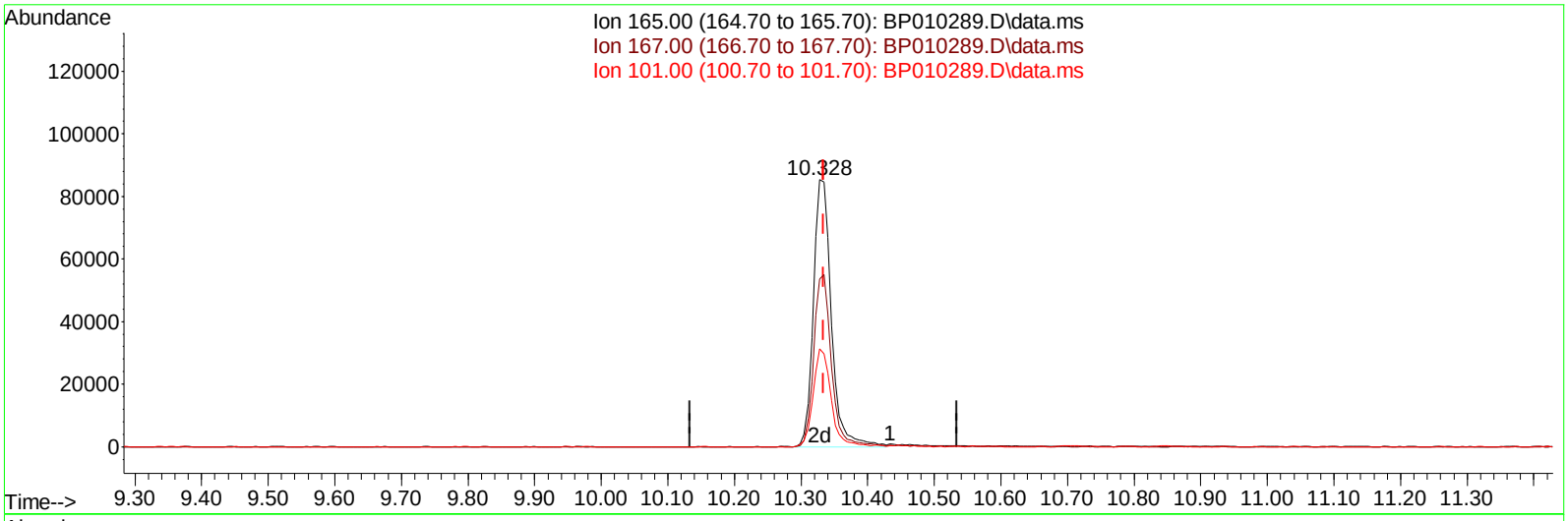

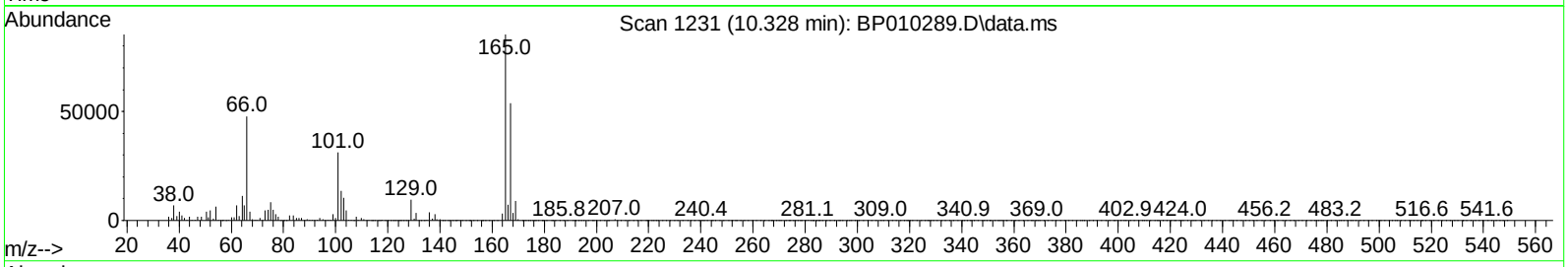

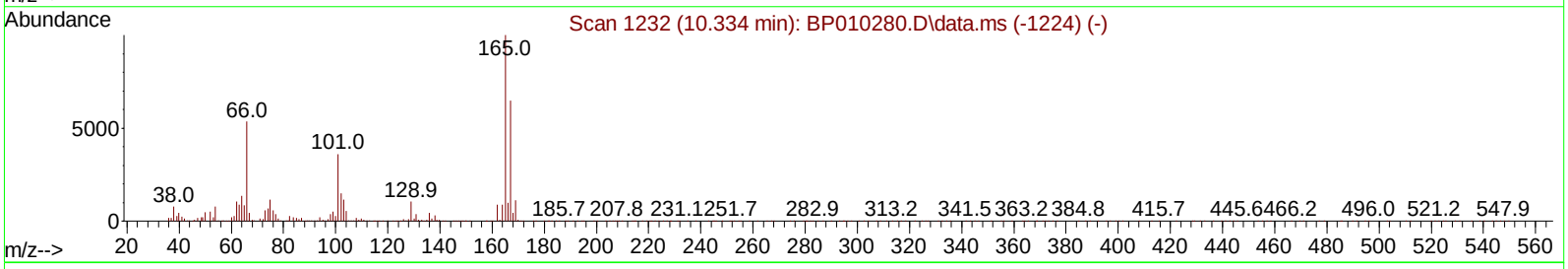

TIC: BP010289.D\data.ms

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

10.328min (-0.006) 16.26 ng/ul m

response 160007

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	62.93
101.00	49.80	36.63#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	131388	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	591337	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	429333	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	948953	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.375	240	927192	20.000	ng/ul	# 0.00
88) Perylene-d12	23.804	264	934747	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	11101m	3.589	ng/uL	0.00
4) Pyridine-d5	3.781	84	36747	4.270	ng/ul	0.01
7) Phenol-d5	7.069	99	194638	15.950	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	141638	18.762	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	159980	17.801	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	164671	16.414	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	86048	18.148	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	91734	16.969	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	160007m	16.262	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	232022	15.973	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	679395	19.863	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	713821	17.504	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	79007	12.042	ng/ul	0.00
60) Fluorene-d10	15.528	176	567573	19.658	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	87190	13.417	ng/ul	0.00
73) Anthracene-d10	17.386	188	941019	20.250	ng/ul	0.00
81) Pyrene-d10	19.622	212	1219375	23.799	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.651	264	1143257	22.934	ng/ul	0.00
Target Compounds						Qvalue

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

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