Quantitation Report (Qedit)

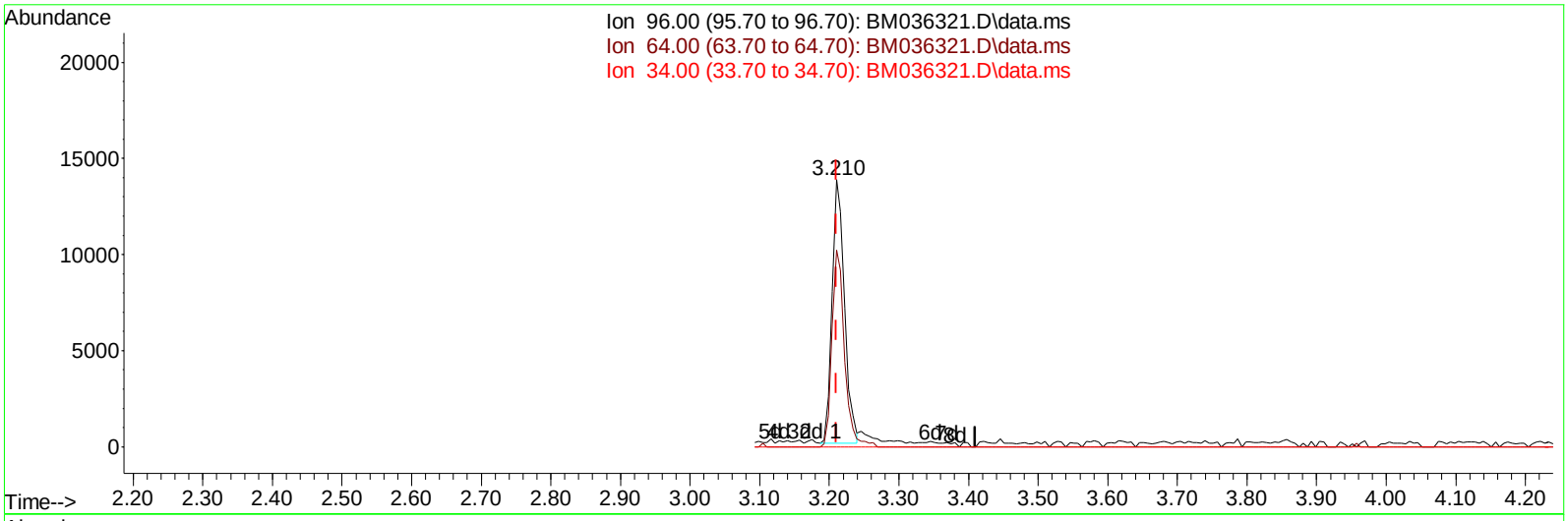
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081722\
Data File : BM036321.D
Acq On : 17 Aug 2022 18:00
Operator : CG/JU
Sample : N4173-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

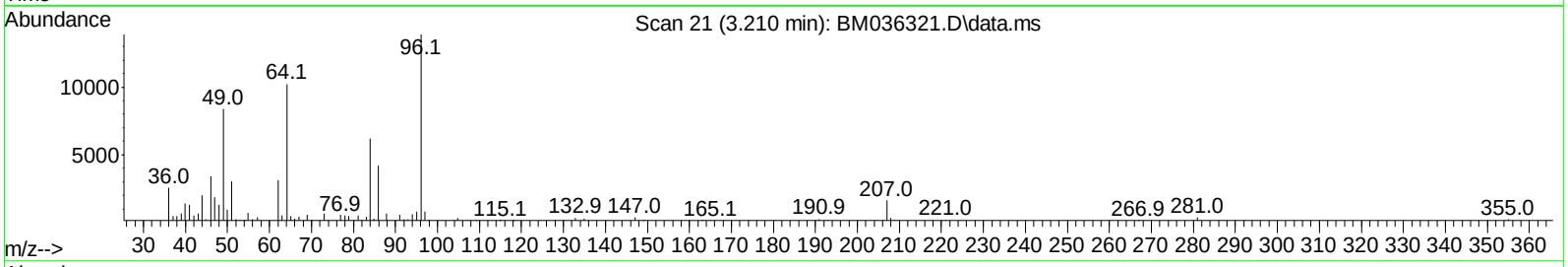
Quant Time: Aug 18 02:18:38 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 16 15:00:33 2022
Response via : Initial Calibration

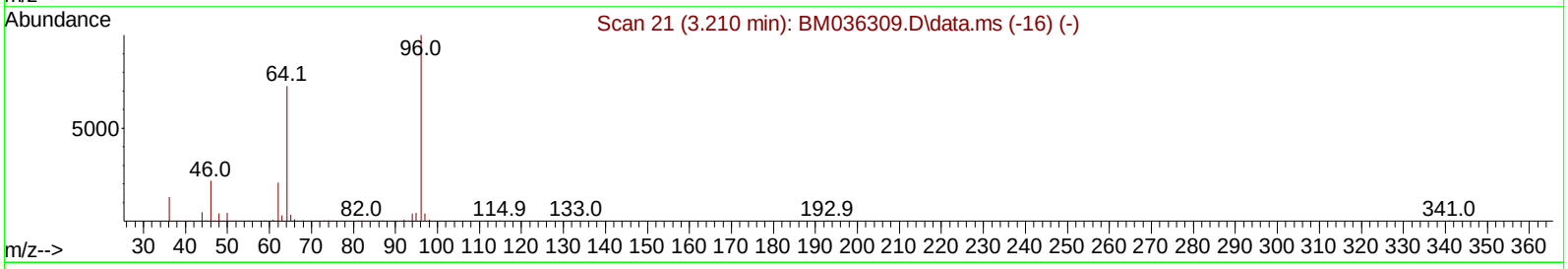
Instrument :
BNA_M
ClientSampleId :
EX7N9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/18/2022
Supervised By :Sohil Jodhani 08/19/2022







TIC: BM036321.D\data.ms

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

3.210min (+ 0.000) 2.70 ng/uL

response 16956

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	73.57
34.00	0.00	0.00
0.00	0.00	0.00

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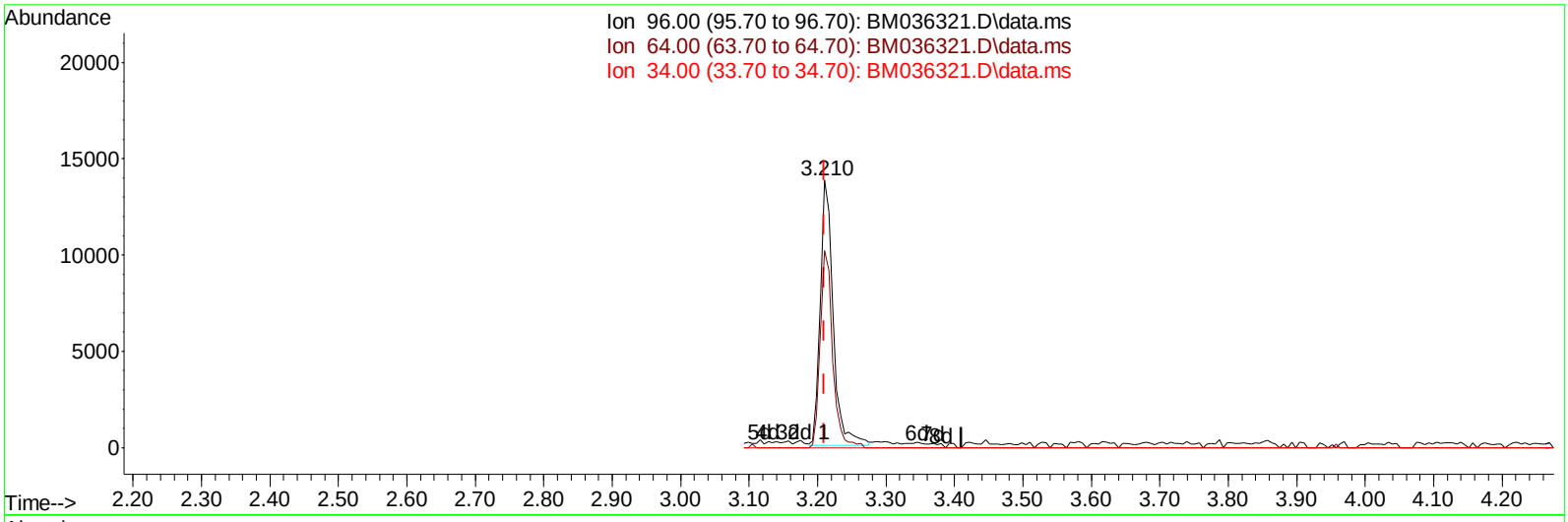

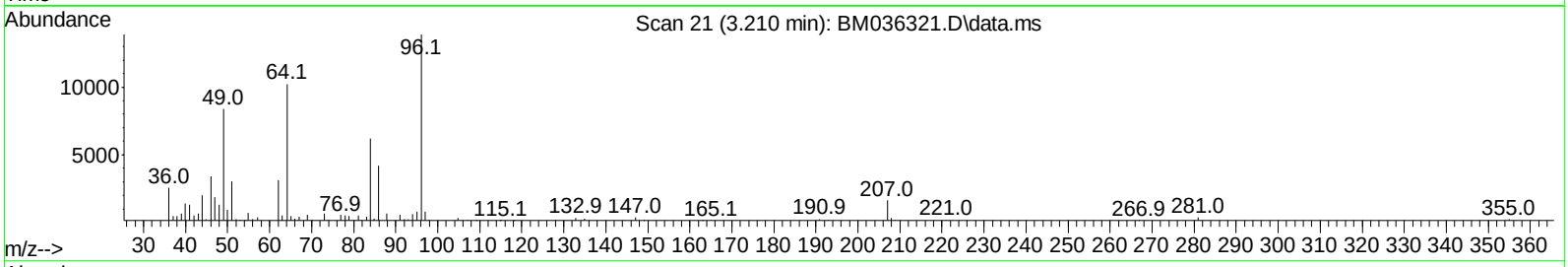

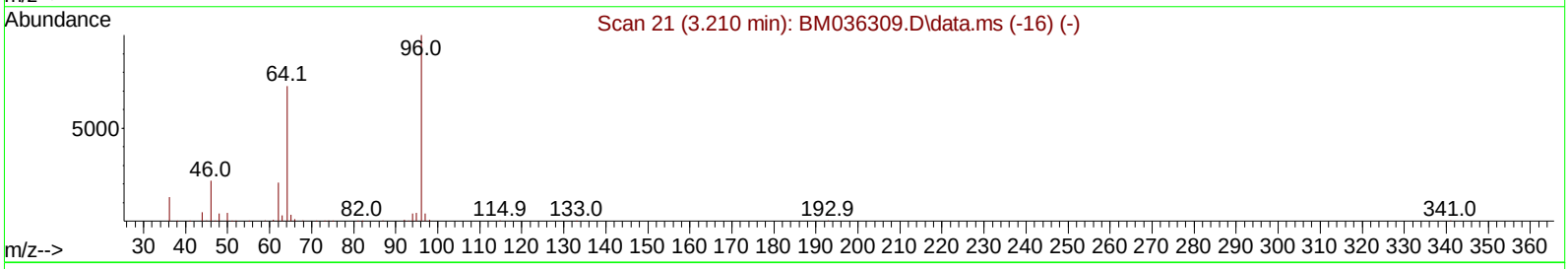

TIC: BM036321.D\data.ms

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

3.210min (+ 0.000) 2.88 ng/uL m

response 18096

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	73.57
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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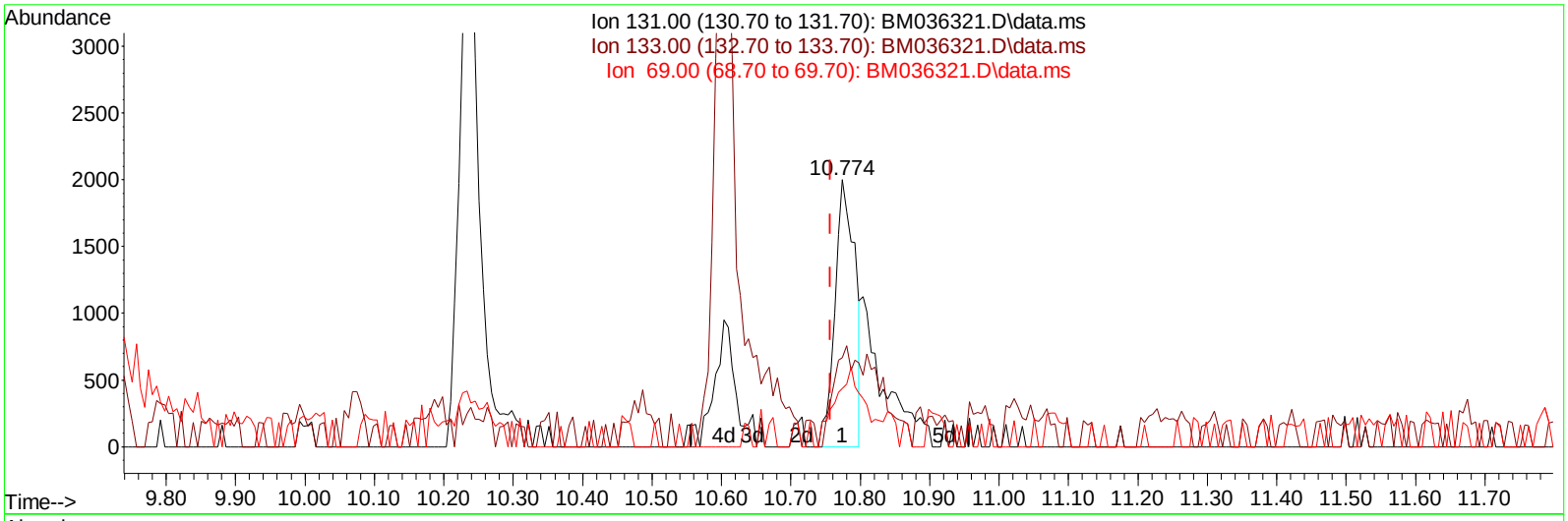

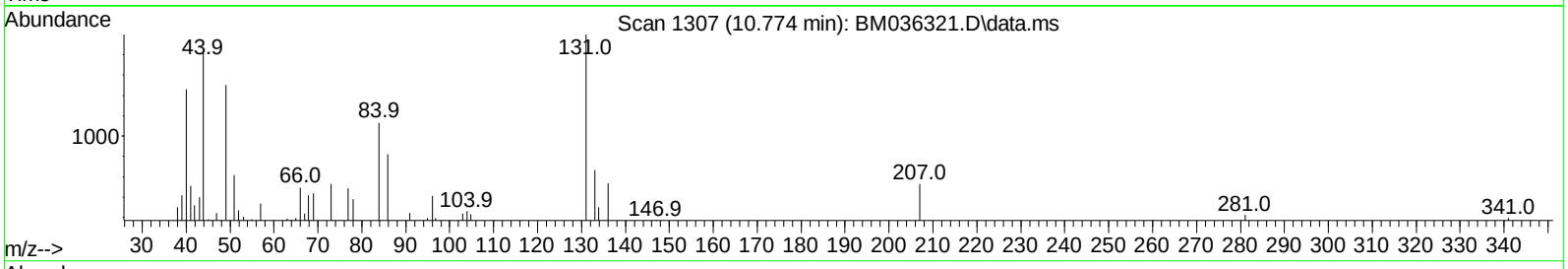

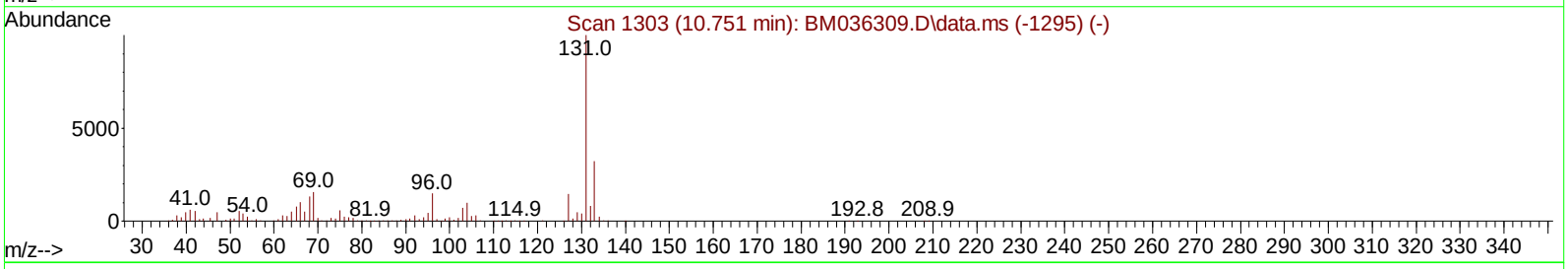

TIC: BM036321.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.774min (+ 0.018) 0.18 ng/ul

response 4037

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.30	33.35
69.00	15.70	21.75#
0.00	0.00	0.00

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ClientSampleId :
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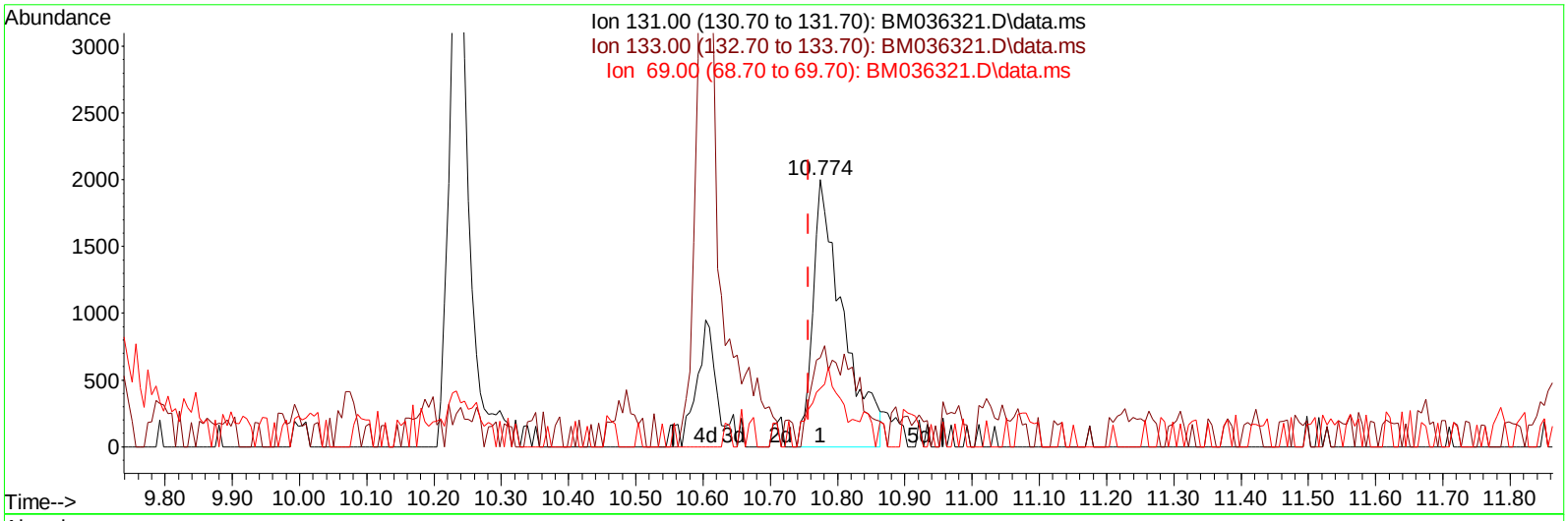

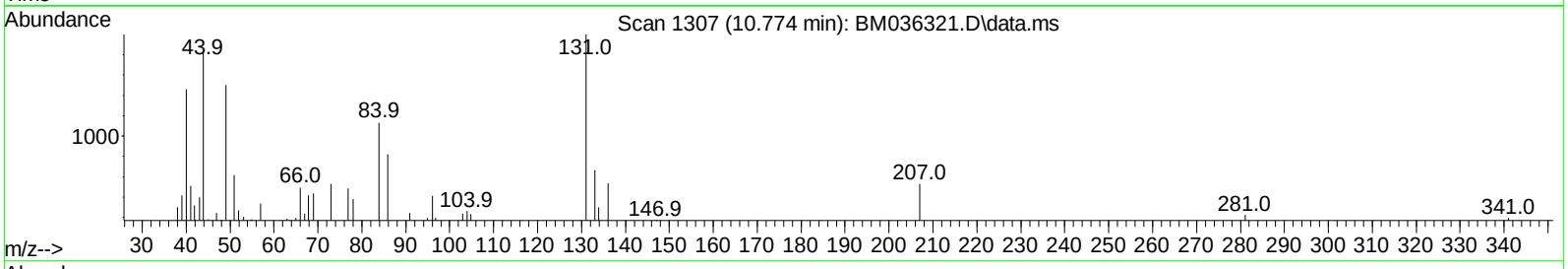

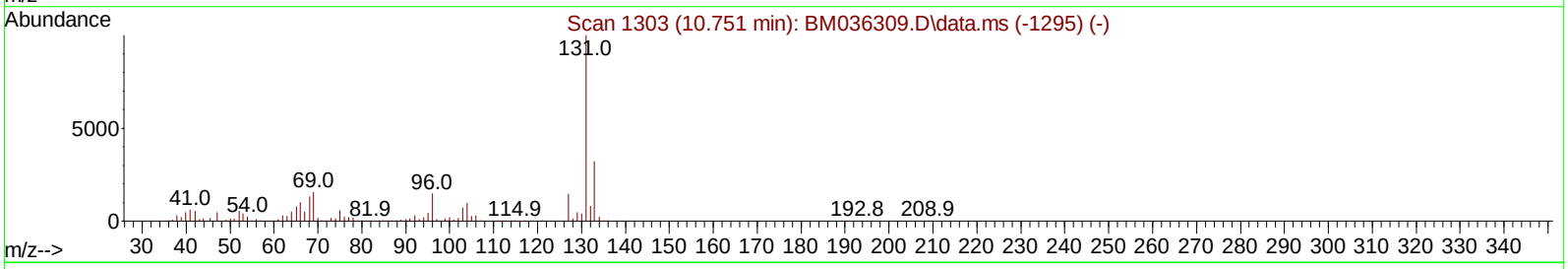

TIC: BM036321.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.774min (+ 0.018) 0.28 ng/ul m

response 6198

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.30	33.35
69.00	15.70	21.75#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/18/2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	239078	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	1055195	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	589951	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	1073690	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	942131	20.000	ng/ul	0.00
88) Perylene-d12	23.703	264	980626	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	18096m	2.885	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.981	99	356074	17.448	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.139	67	236727	18.429	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	297375	18.546	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	272895	17.910	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	147614	18.451	ng/ul	0.00
24) 2-Nitrophenol-d4	9.692	143	150575	18.555	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.233	165	249247	17.189	ng/ul	0.00
31) 4-Chloroaniline-d4	10.774	131	6198m	0.276	ng/ul	0.02
46) Dimethylphthalate-d6	13.863	166	833459	22.331	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	934811	20.082	ng/ul	0.00
54) 4-Nitrophenol-d4	14.680	143	109027	15.417	ng/ul	0.00
60) Fluorene-d10	15.439	176	732203	21.710	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.568	200	85287	15.451	ng/ul	0.00
73) Anthracene-d10	17.286	188	1143681	23.956	ng/ul	0.00
81) Pyrene-d10	19.580	212	1405085	28.750	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.550	264	1354347	28.169	ng/ul	0.00

Target Compounds Qvalue

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

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