Quantitation Report (Qedit)

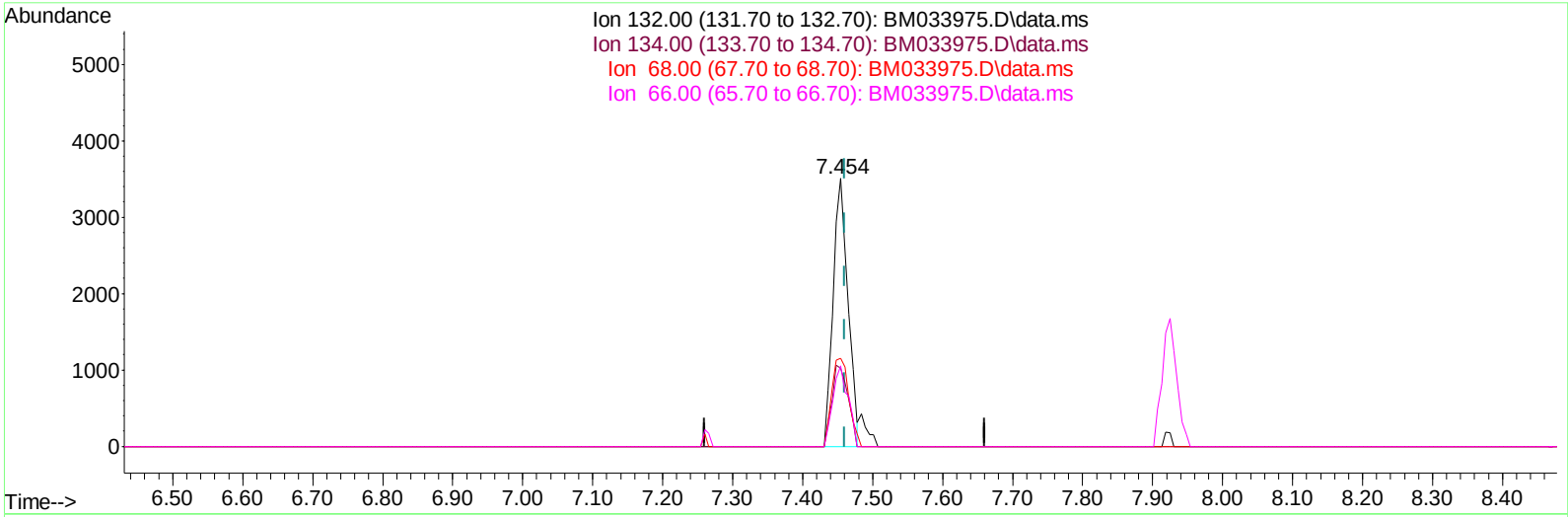
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033975.D
Acq On : 02 Feb 2022 15:19
Operator : CG/JU
Sample : N1307-05DL 30X
Misc :
ALS Vial : 46 Sample Multiplier: 1

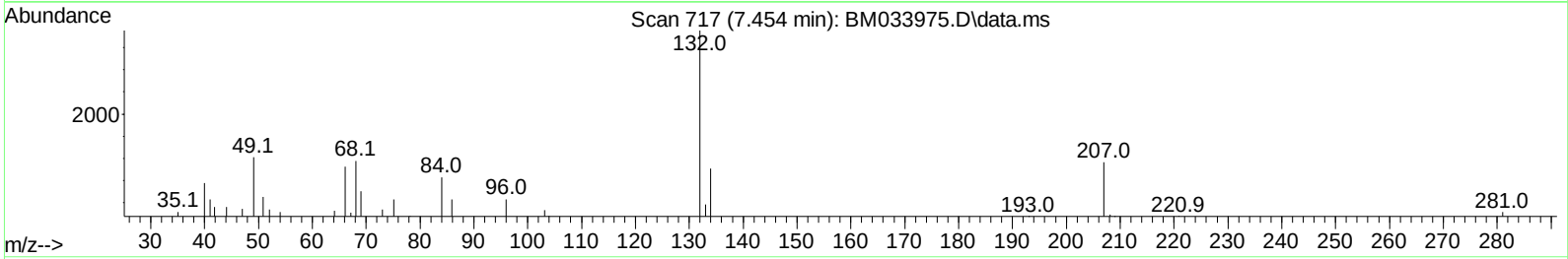
Quant Time: Feb 03 00:13:14 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

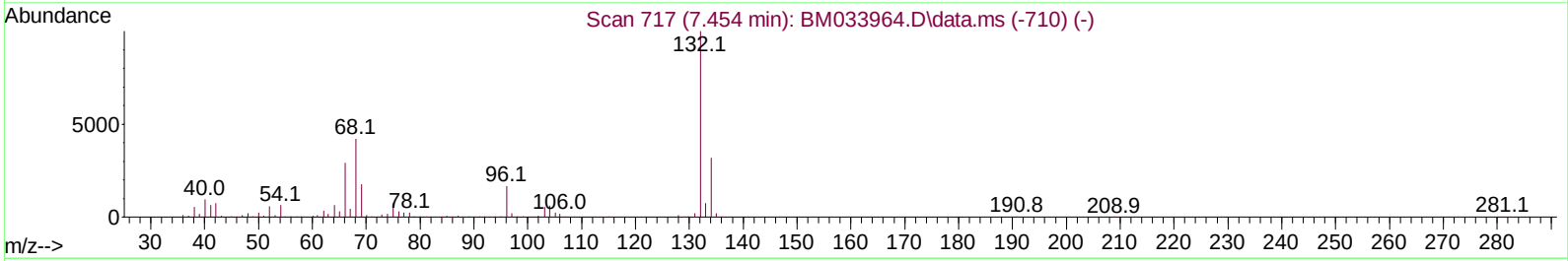
Instrument :
BNA_M
ClientSampleId :
C0Q38DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :Yogesh Patel 02/04/2022







TIC: BM033975.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.454min (-0.006) 0.75 ng/ul

response 5110

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	30.30	29.03
68.00	48.30	32.91#
66.00	37.00	30.08

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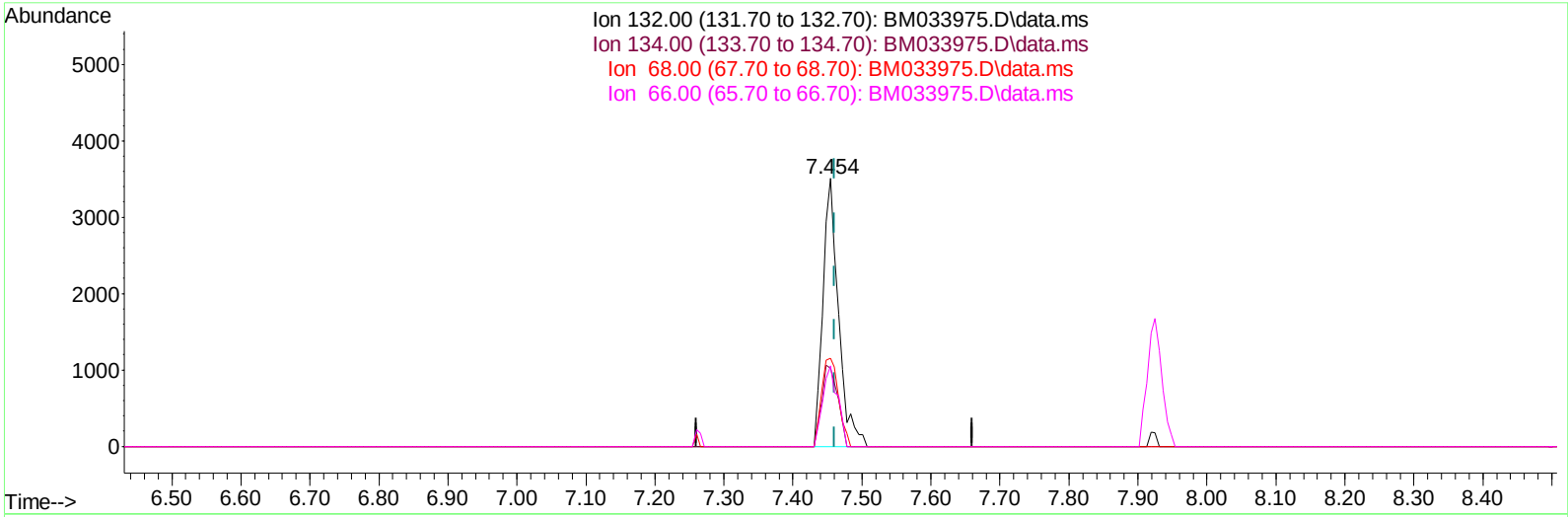

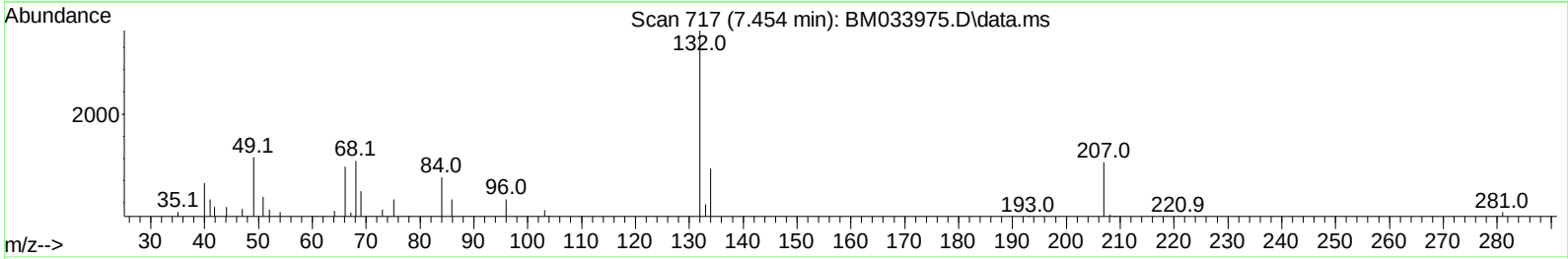

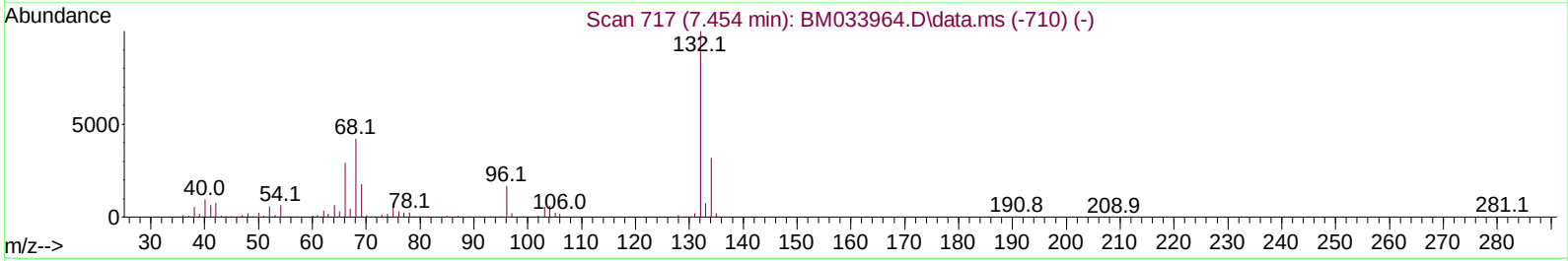

TIC: BM033975.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.454min (-0.006) 0.80 ng/ul m

response 5456

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	30.30	29.03
68.00	48.30	32.91#
66.00	37.00	30.08

Quantitation Report (Qedit)

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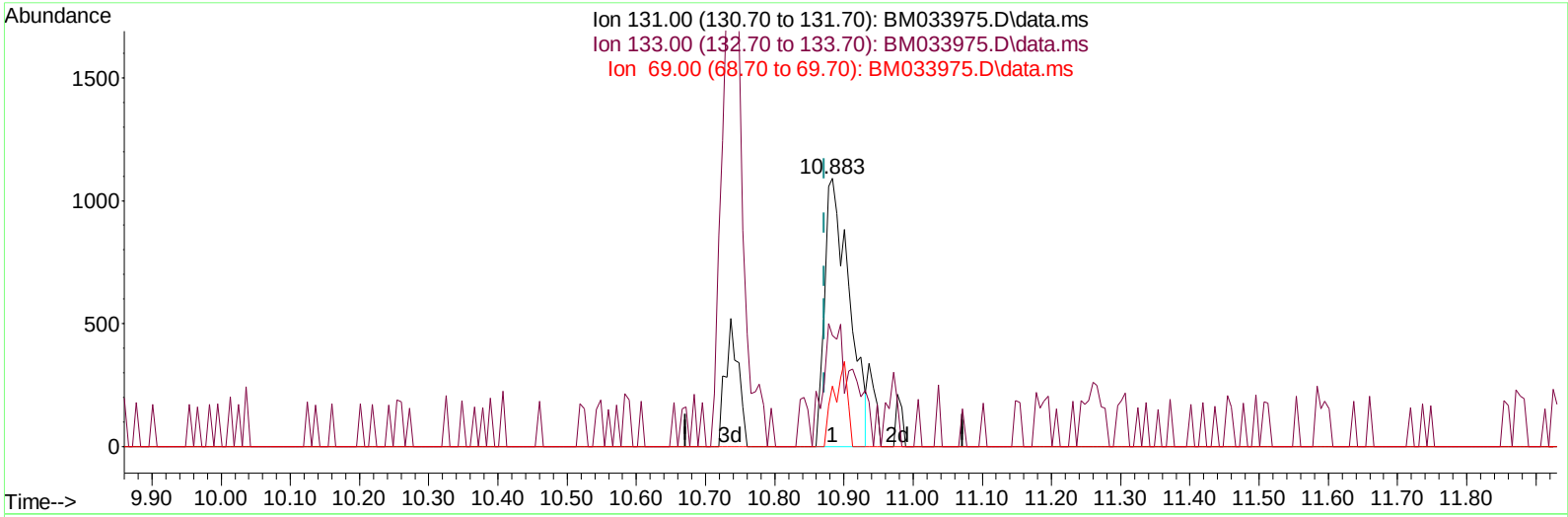

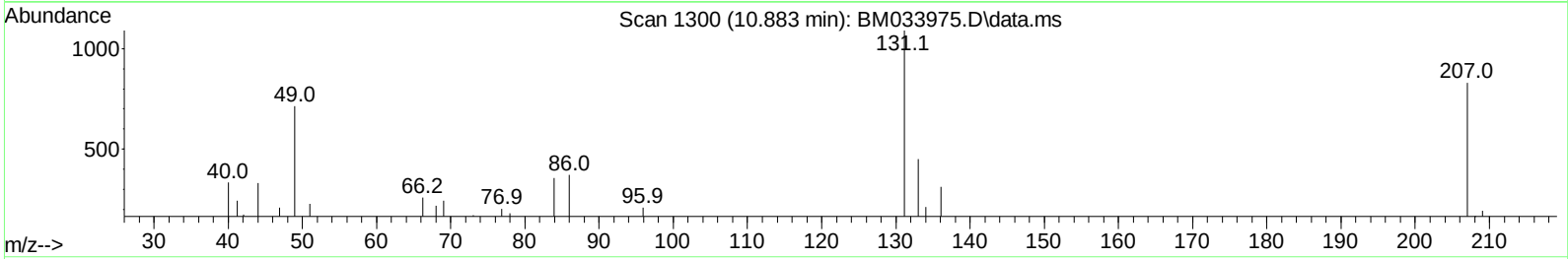

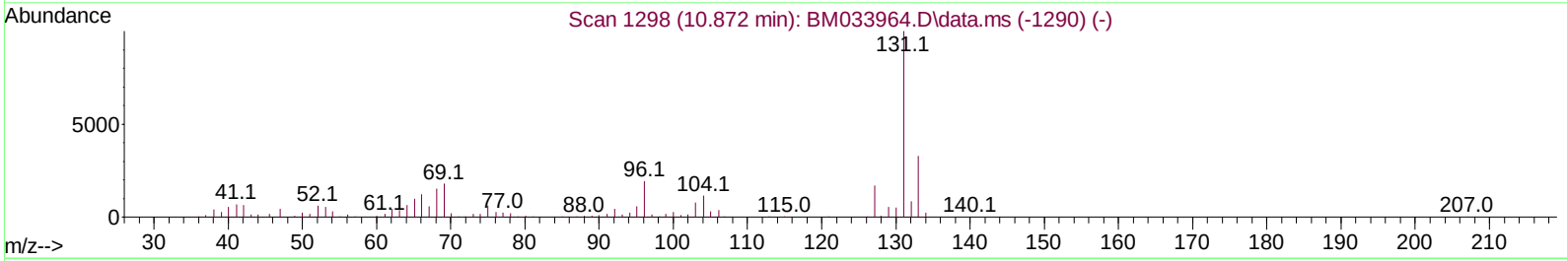

TIC: BM033975.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.883min (+ 0.012) 0.28 ng/ul

response 2679

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.10	41.38#
69.00	22.50	22.48
0.00	0.00	0.00

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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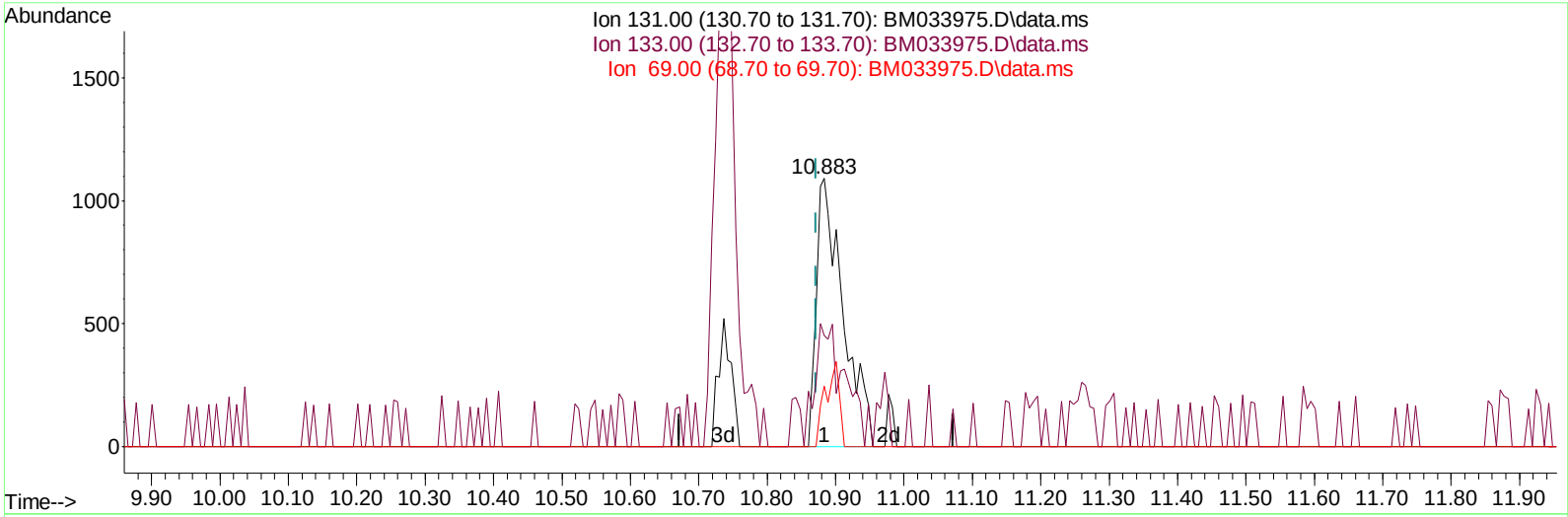

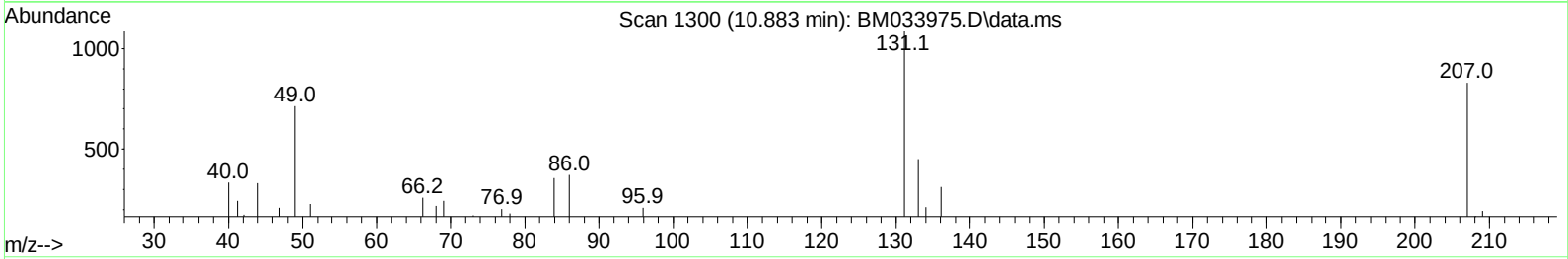

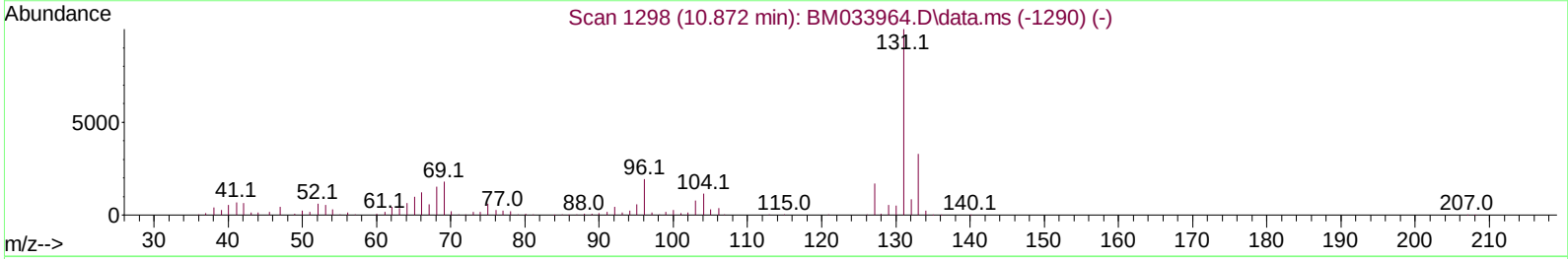

TIC: BM033975.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.883min (+ 0.012) 0.30 ng/ul m

response 2943

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.10	41.38#
69.00	22.50	22.48
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	101186	20.000	ng/ul	0.00
20) Naphthalene-d8	10.736	136	449239	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.560	164	326084	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	694060	20.000	ng/ul	0.00
79) Chrysene-d12	21.459	240	546893	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	477293	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.090	99	1367	0.161	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	5839	1.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	5456m	0.803	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	2622	0.390	ng/ul	0.00
21) Nitrobenzene-d5	9.089	128	2788	0.773	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	2399	0.616	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.372	165	4252	0.604	ng/ul	0.01
31) 4-Chloroaniline-d4	10.883	131	2943m	0.303	ng/ul	0.01
46) Dimethylphthalate-d6	13.971	166	29576	1.214	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	28793	0.965	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.548	176	23494	1.155	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	3072	0.705	ng/ul	0.00
73) Anthracene-d10	17.395	188	43251	1.301	ng/ul	0.00
81) Pyrene-d10	19.677	212	40910	1.267	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.694	264	27428	1.094	ng/ul	-0.01
Target Compounds						Qvalue
58) 2,3,4,6-Tetrachlorophenol	15.189	232	6808	1.208	ng/ul#	94
71) Pentachlorophenol	16.948	266	327312	59.514	ng/ul	96

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

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