Quantitation Report (Qedit)

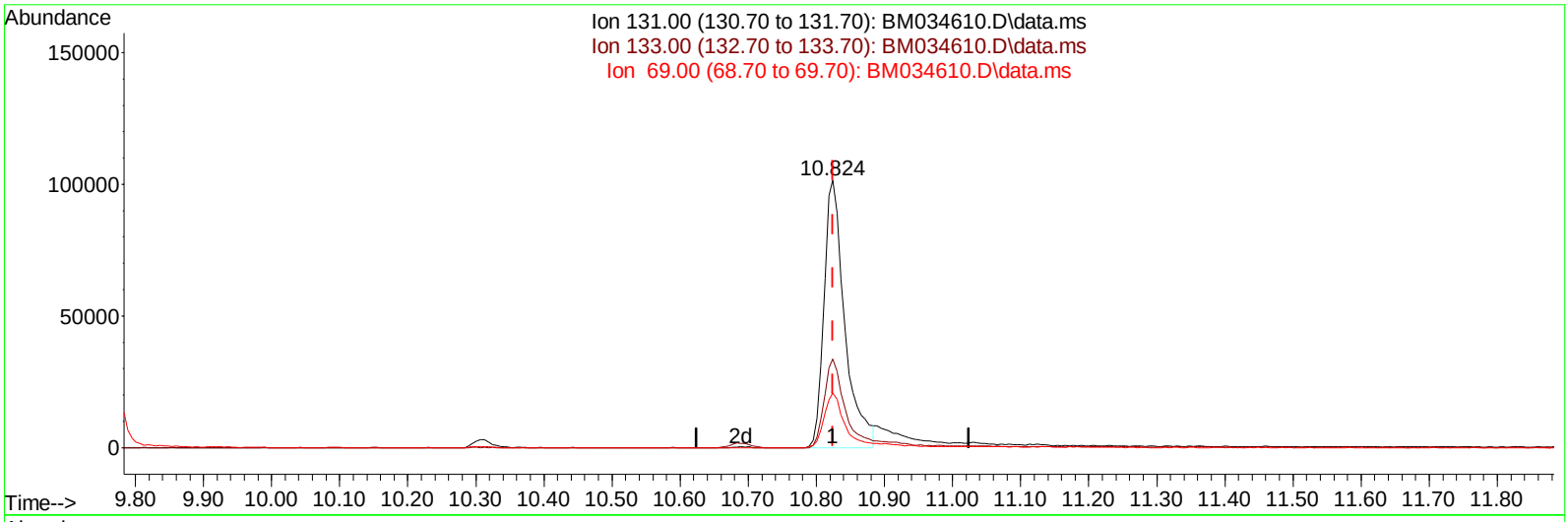
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034610.D
Acq On : 11 Apr 2022 15:15
Operator : CG/JU
Sample : N2338-06
Misc :
ALS Vial : 8 Sample Multiplier: 1

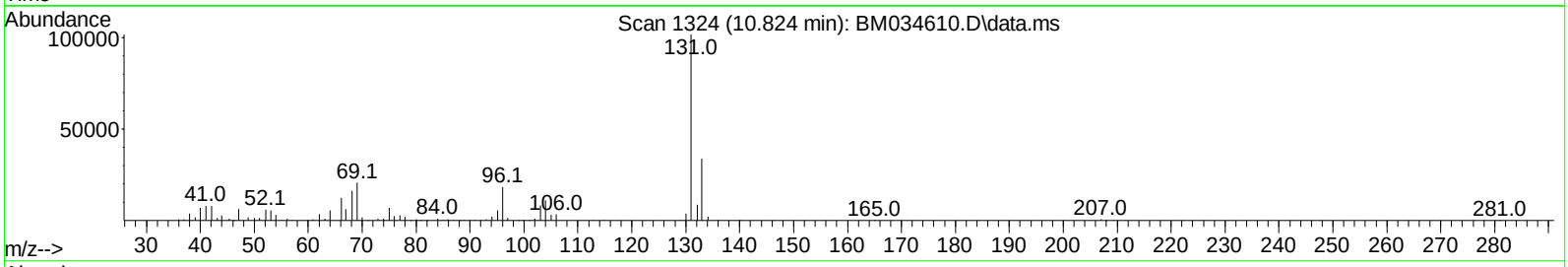
Quant Time: Apr 12 02:55:17 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 08 14:30:08 2022
Response via : Initial Calibration

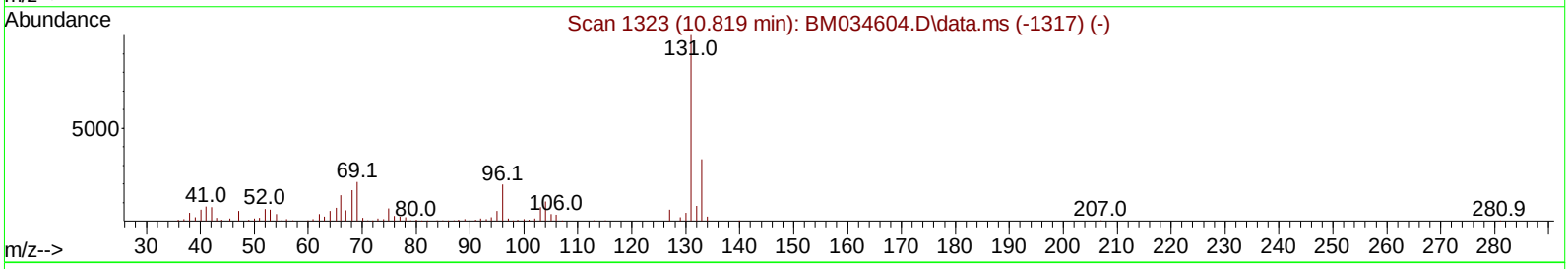
Instrument :
BNA_M
ClientSampleId :
C0J91

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/14/2022







TIC: BM034610.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.824min (-0.000) 27.56 ng/ul

response 214703

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.18
69.00	16.80	20.33#
0.00	0.00	0.00

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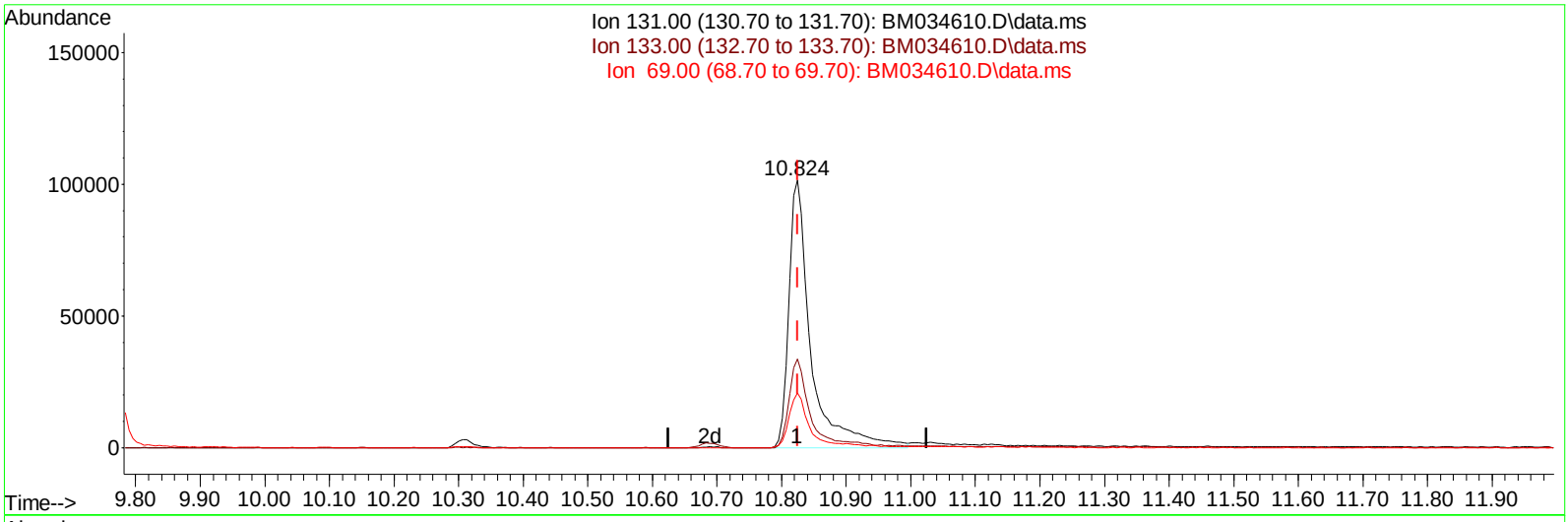

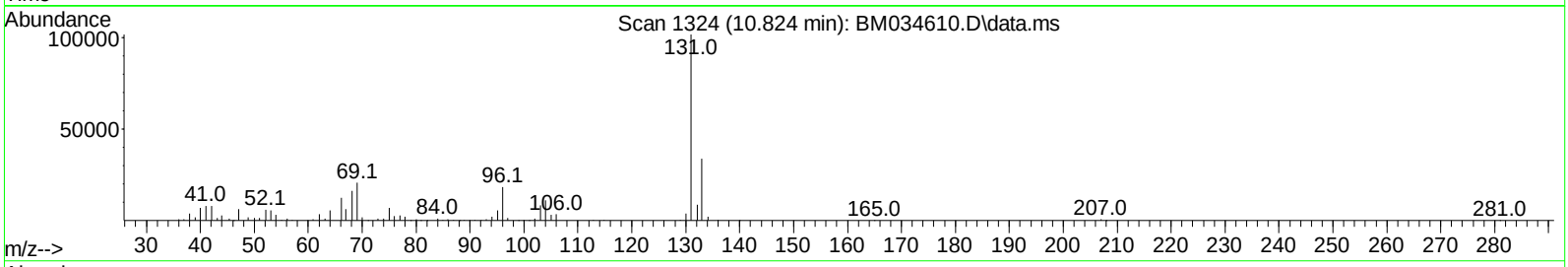

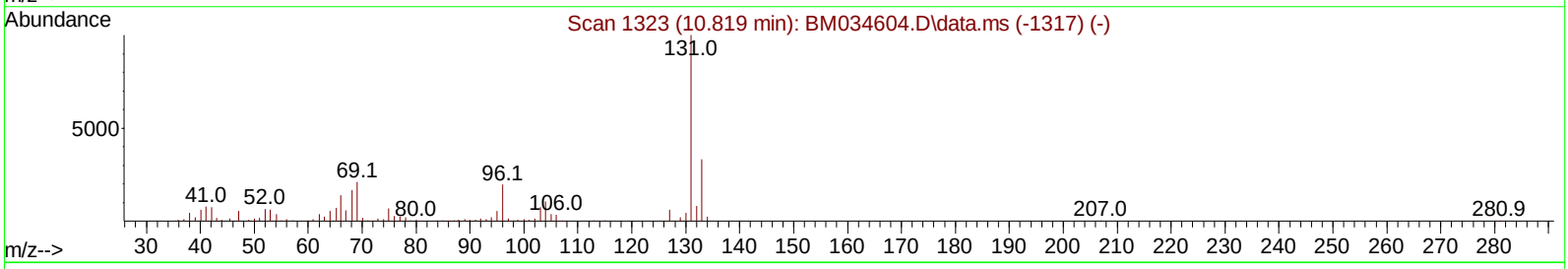

TIC: BM034610.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.824min (-0.000) 31.04 ng/ul m

response 241820

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.18
69.00	16.80	20.33#
0.00	0.00	0.00

Quantitation Report (Qedit)

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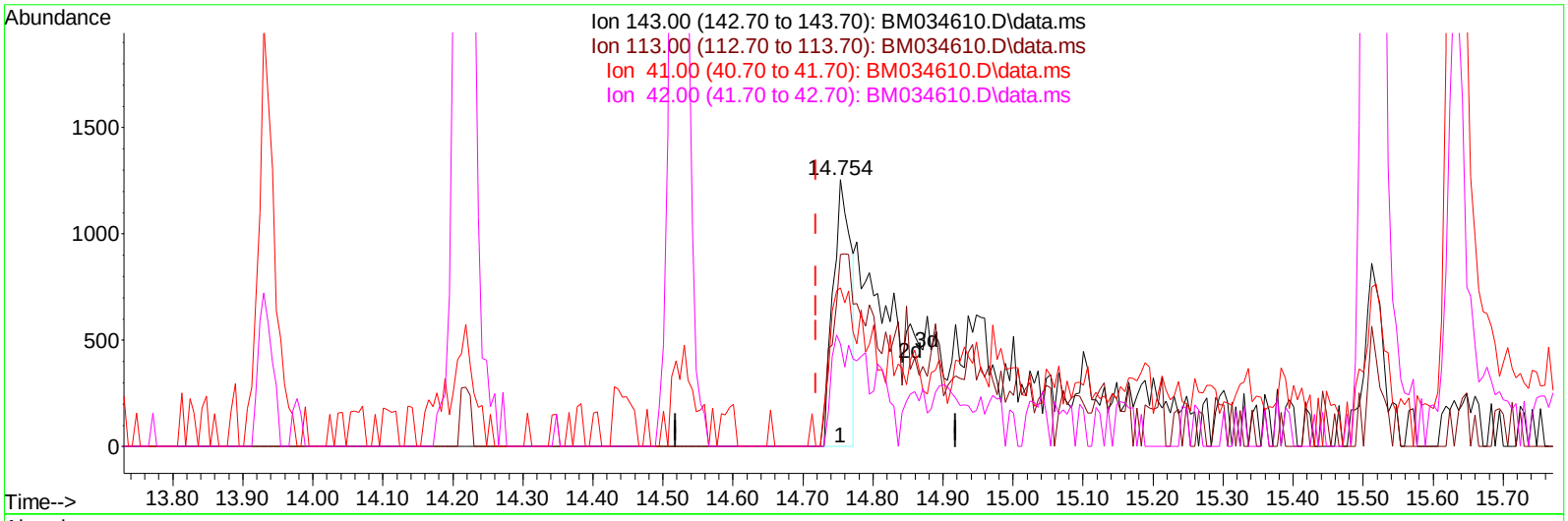

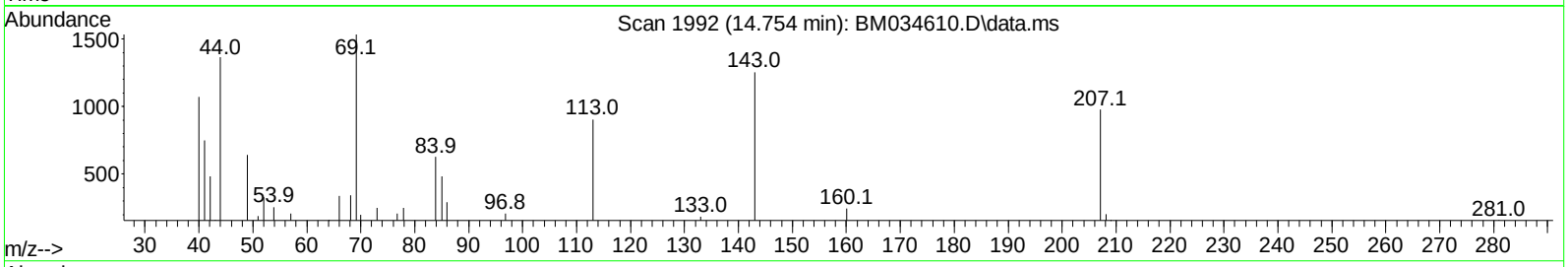

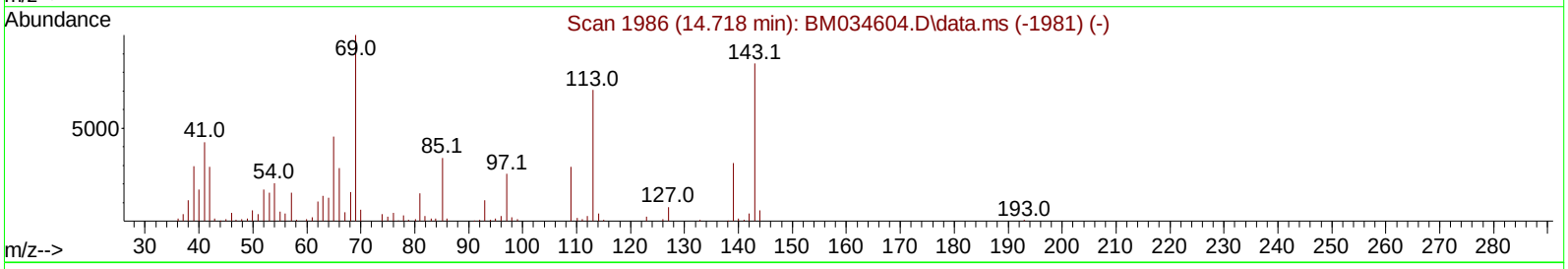

TIC: BM034610.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.035) 0.89 ng/ul

response 2236

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	72.11
41.00	31.90	59.60#
42.00	19.10	38.57#

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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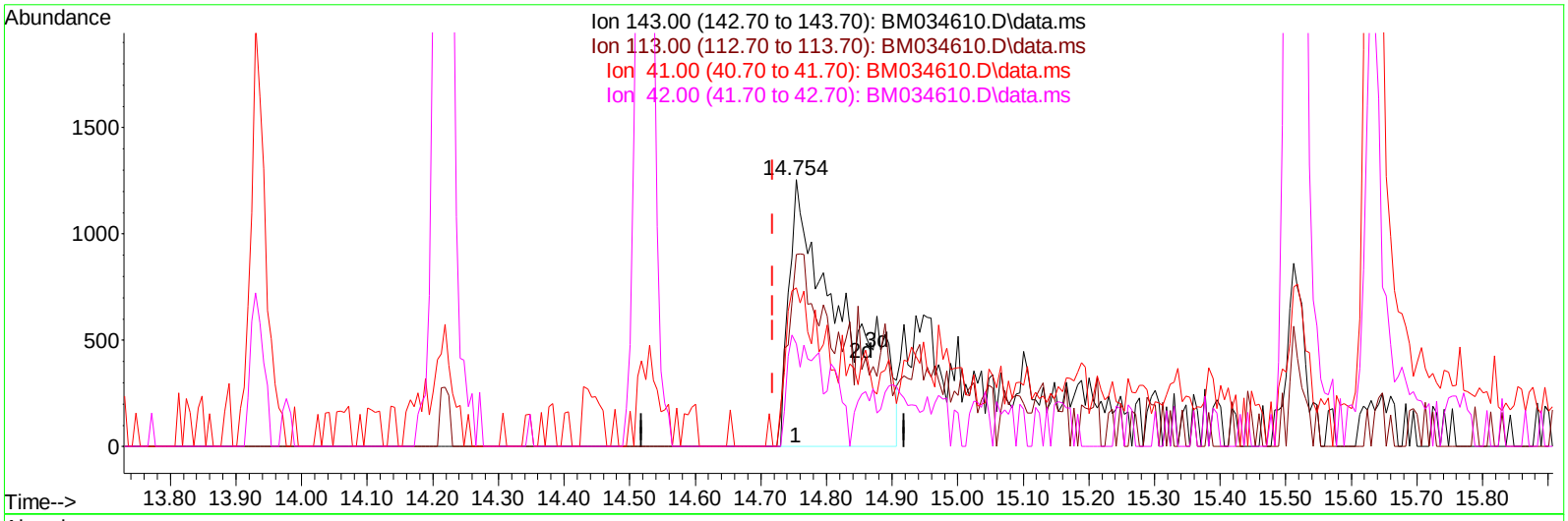

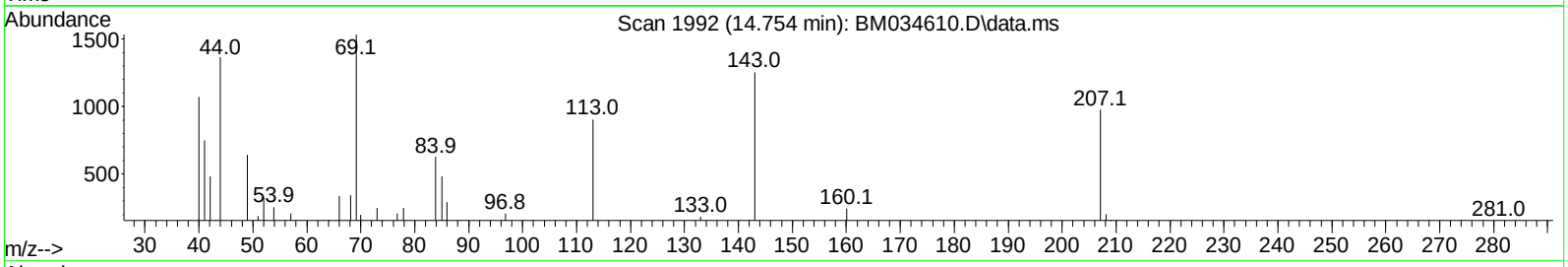

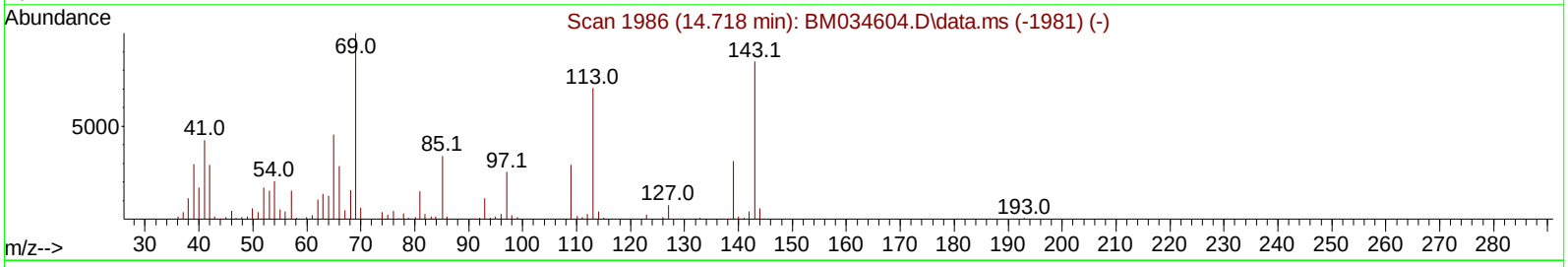

TIC: BM034610.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.035) 2.78 ng/ul m

response 7011

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	72.11
41.00	31.90	59.60#
42.00	19.10	38.57#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 04/14/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	95830	20.000	ng/ul	0.00
20) Naphthalene-d8	10.689	136	391889	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.524	164	229247	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	450768	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.441	240	442190	20.000	ng/ul	# 0.00
88) Perylene-d12	23.824	264	492742	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.254	96	14549	5.414	ng/uL	0.00
4) Pyridine-d5	3.684	84	77264	11.913	ng/ul	0.00
7) Phenol-d5	7.042	99	61435	7.702	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.201	67	202313	38.065	ng/ul	0.00
11) 2-Chlorophenol-d4	7.407	132	189824	30.787	ng/ul	0.00
15) 4-Methylphenol-d8	8.589	113	115702	18.570	ng/ul	0.00
21) Nitrobenzene-d5	9.042	128	112726	39.139	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	110558	36.870	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	176338	31.230	ng/ul	0.00
31) 4-Chloroaniline-d4	10.824	131	241820m	31.043	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	681003	41.111	ng/ul	0.00
49) Acenaphthylene-d8	14.218	160	812034	38.140	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	7011m	2.776	ng/ul	0.04
60) Fluorene-d10	15.512	176	594284	41.515	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.636	200	86802	31.543	ng/ul	0.00
73) Anthracene-d10	17.365	188	950731	44.238	ng/ul	0.00
81) Pyrene-d10	19.653	212	1169826	48.471	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.671	264	1085527	43.796	ng/ul	0.00

Target Compounds Qvalue

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

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