Quantitation Report (Qedit)

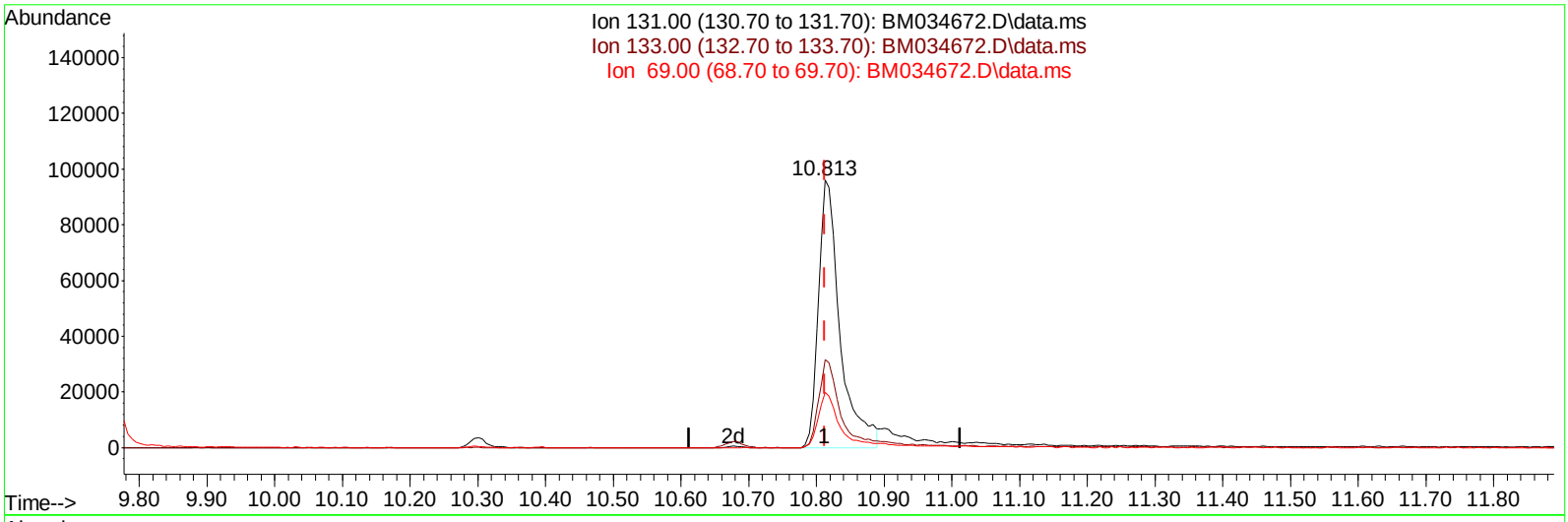
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034672.D
Acq On : 14 Apr 2022 17:21
Operator : CG/JU
Sample : N2266-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

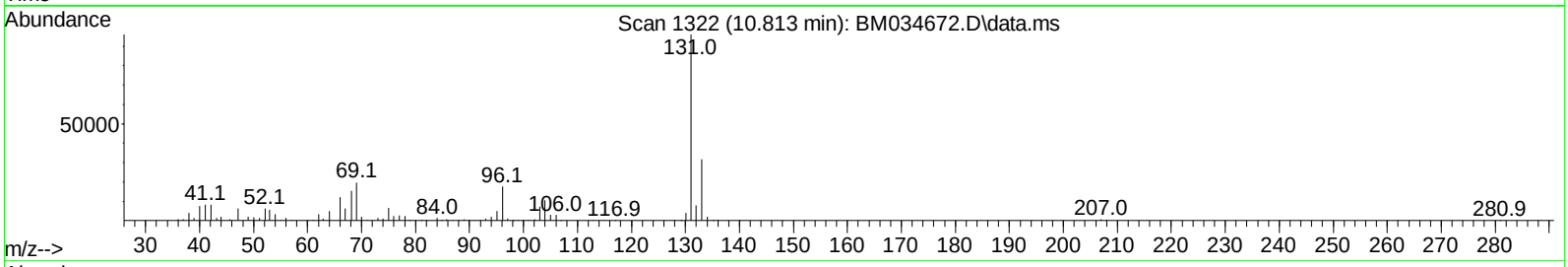
Quant Time: Apr 15 00:47:01 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 14 23:42:00 2022
Response via : Initial Calibration

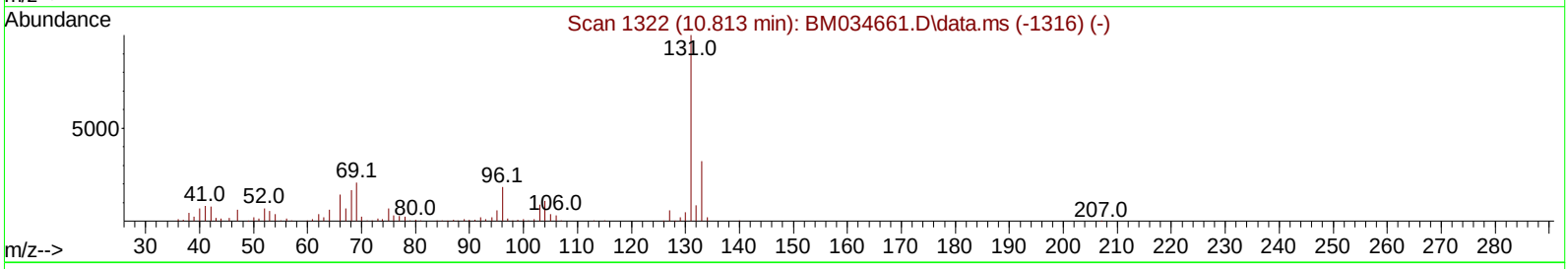
Instrument :
BNA_M
ClientSampleId :
YAZ04

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/23/2022







TIC: BM034672.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.813min (-0.000) 22.75 ng/ul

response 212973

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.03
69.00	16.80	20.64#
0.00	0.00	0.00

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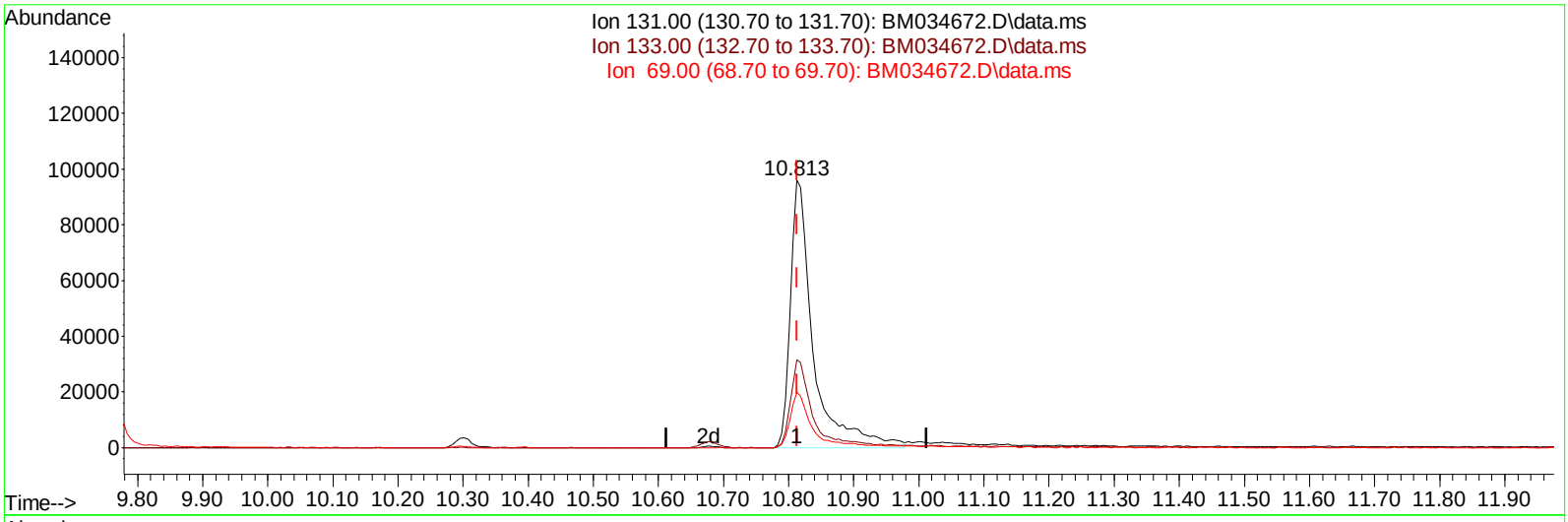

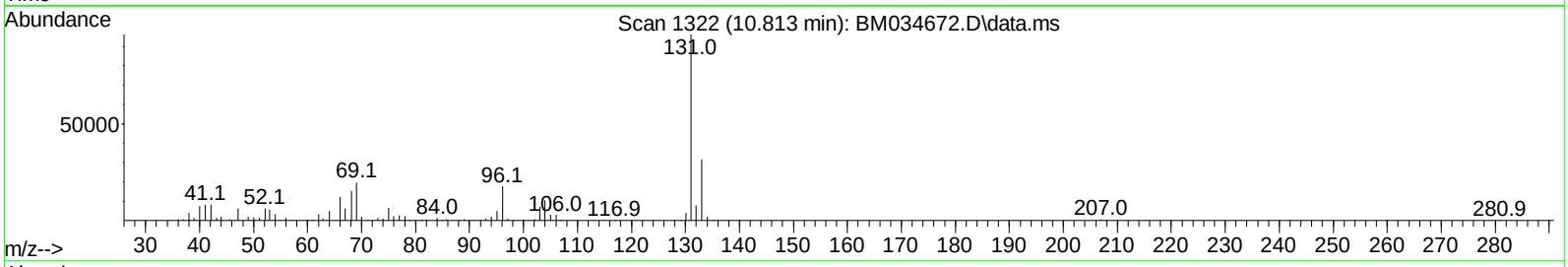

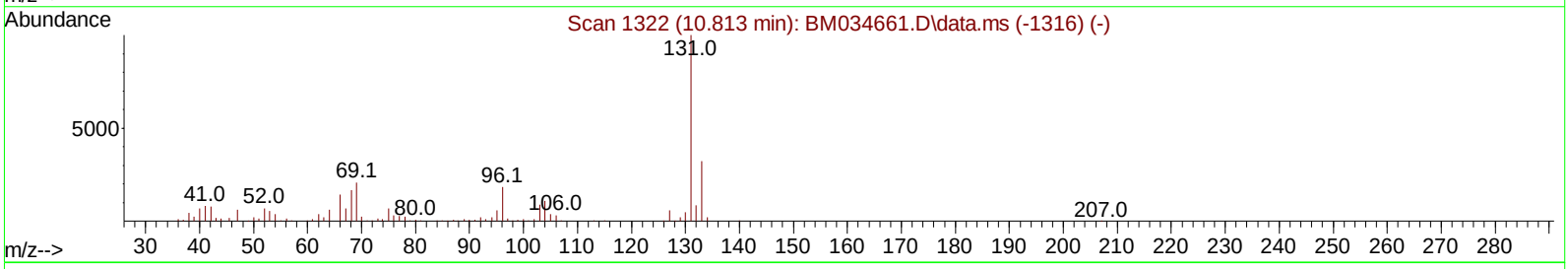

TIC: BM034672.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.813min (-0.000) 24.98 ng/ul m

response 233871

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	33.03
69.00	16.80	20.64#
0.00	0.00	0.00

Quantitation Report (Qedit)

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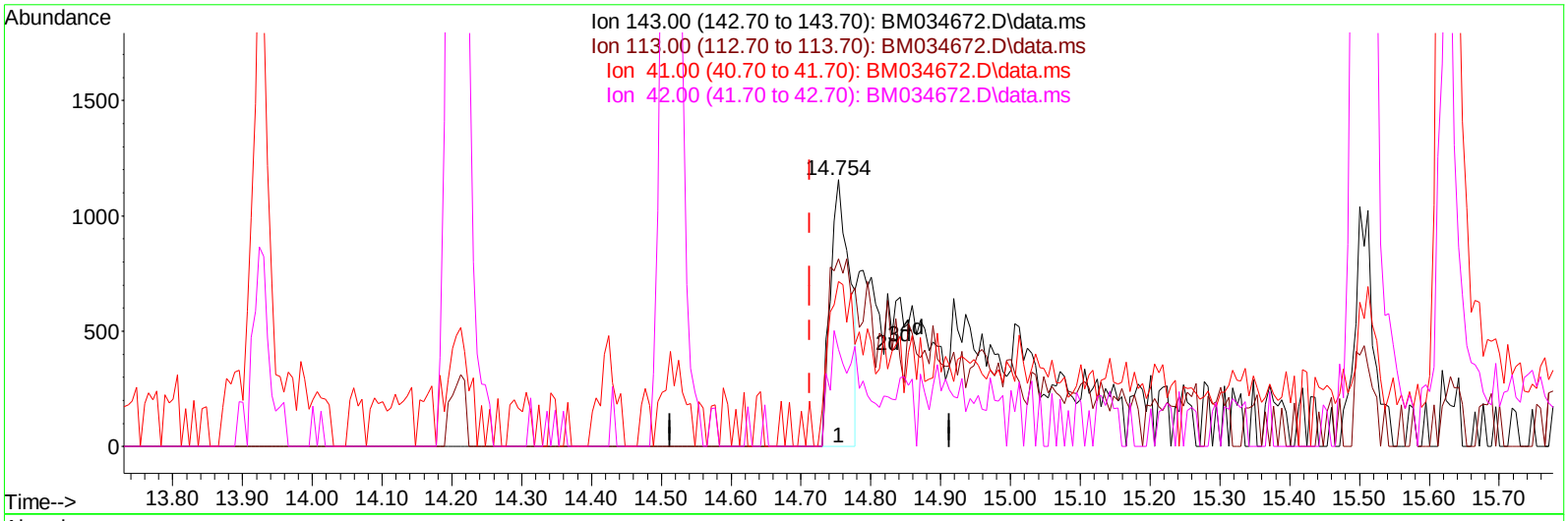

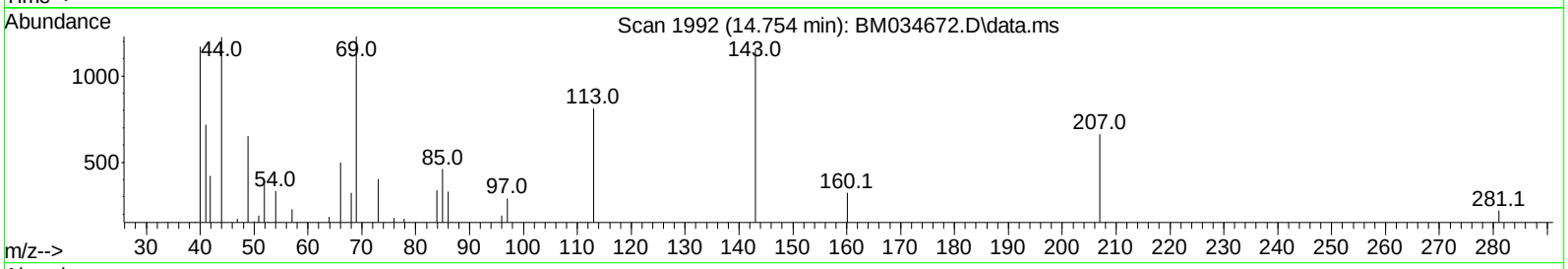

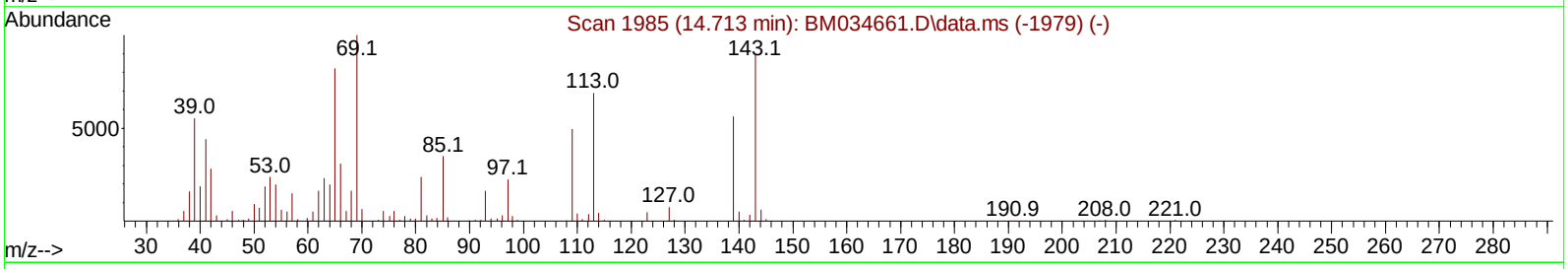

TIC: BM034672.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.041) 0.75 ng/ul

response 2255

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	70.35
41.00	31.90	61.88#
42.00	19.10	36.39#

Quantitation Report (Qedit)

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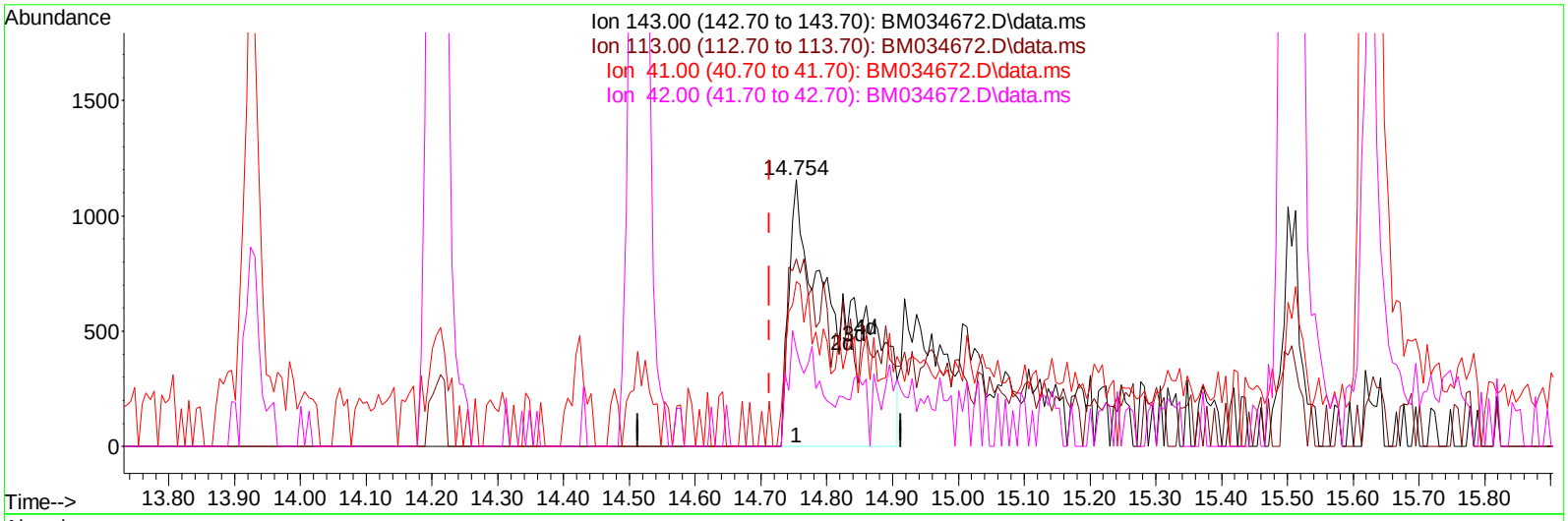

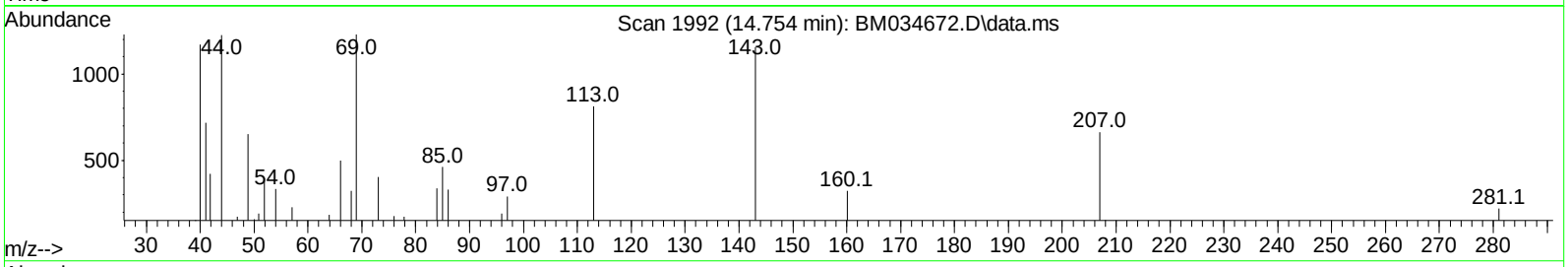

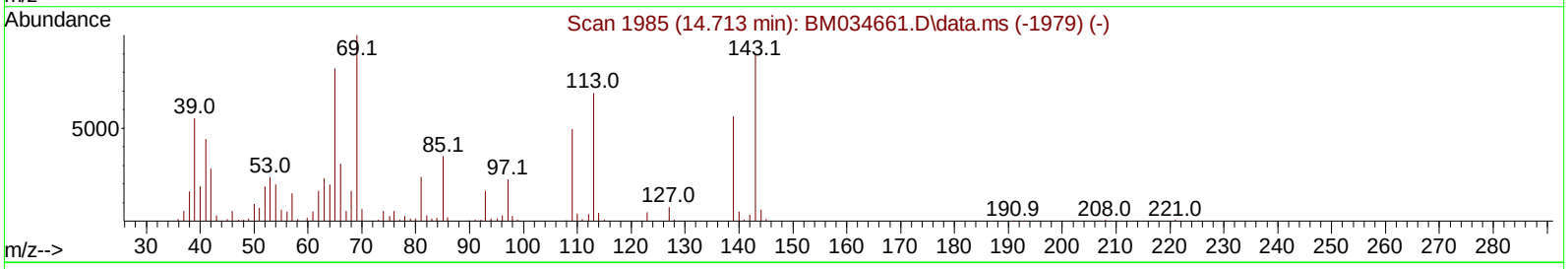

TIC: BM034672.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.041) 2.18 ng/ul m

response 6565

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	70.35
41.00	31.90	61.88#
42.00	19.10	36.39#

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	7.866	152	114608	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.678	136	471002	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.518	164	273387	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.254	188	531710	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.436	240	488913	20.000	ng/ul	# 0.00
88) Perylene-d12	23.812	264	556971	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.249	96	17854	5.555	ng/uL	0.00
4) Pyridine-d5	3.678	84	57420	7.403	ng/ul	0.00
7) Phenol-d5	7.031	99	63706	6.678	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.195	67	211444	33.265	ng/ul	0.00
11) 2-Chlorophenol-d4	7.395	132	189030	25.635	ng/ul	0.00
15) 4-Methylphenol-d8	8.584	113	126099	16.923	ng/ul	0.00
21) Nitrobenzene-d5	9.031	128	111802	32.298	ng/ul	0.00
24) 2-Nitrophenol-d4	9.760	143	110487	30.657	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.301	165	177028	26.086	ng/ul	0.00
31) 4-Chloroaniline-d4	10.813	131	233871m	24.980	ng/ul	0.00
46) Dimethylphthalate-d6	13.924	166	681308	34.489	ng/ul	0.00
49) Acenaphthylene-d8	14.207	160	801193	31.555	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	6565m	2.180	ng/ul	0.04
60) Fluorene-d10	15.507	176	606851	35.549	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.624	200	92011	28.346	ng/ul	0.00
73) Anthracene-d10	17.354	188	924928	36.486	ng/ul	0.00
81) Pyrene-d10	19.648	212	1134782	42.526	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.659	264	1068969	38.154	ng/ul	0.00
Target Compounds						Qvalue

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

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