Quantitation Report (Qedit)

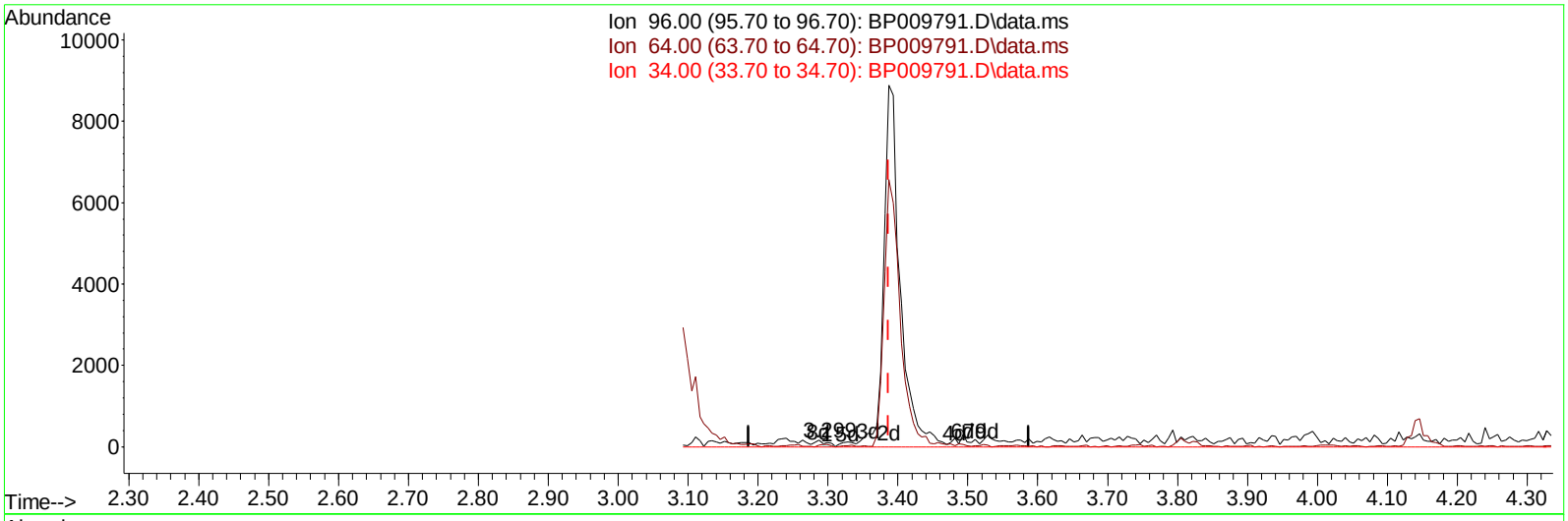
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009791.D
Acq On : 12 Apr 2022 19:50
Operator : CG/JU
Sample : N2342-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

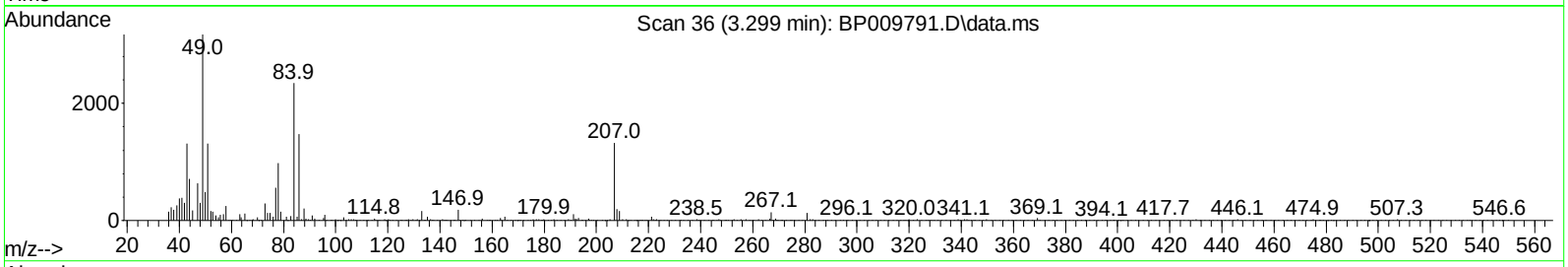
Quant Time: Apr 13 02:25:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

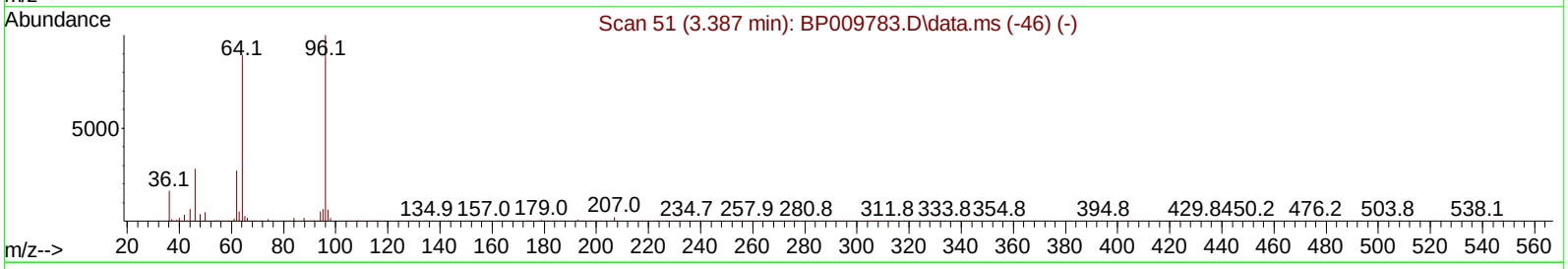
Instrument :
BNA_P
ClientSampleId :
C0J94

Manual IntegrationsAPPROVED

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







TIC: BP009791.D\data.ms

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

3.299min (-0.088) 0.02 ng/uL

response 61

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.91
34.00	0.00	0.00
0.00	0.00	0.00

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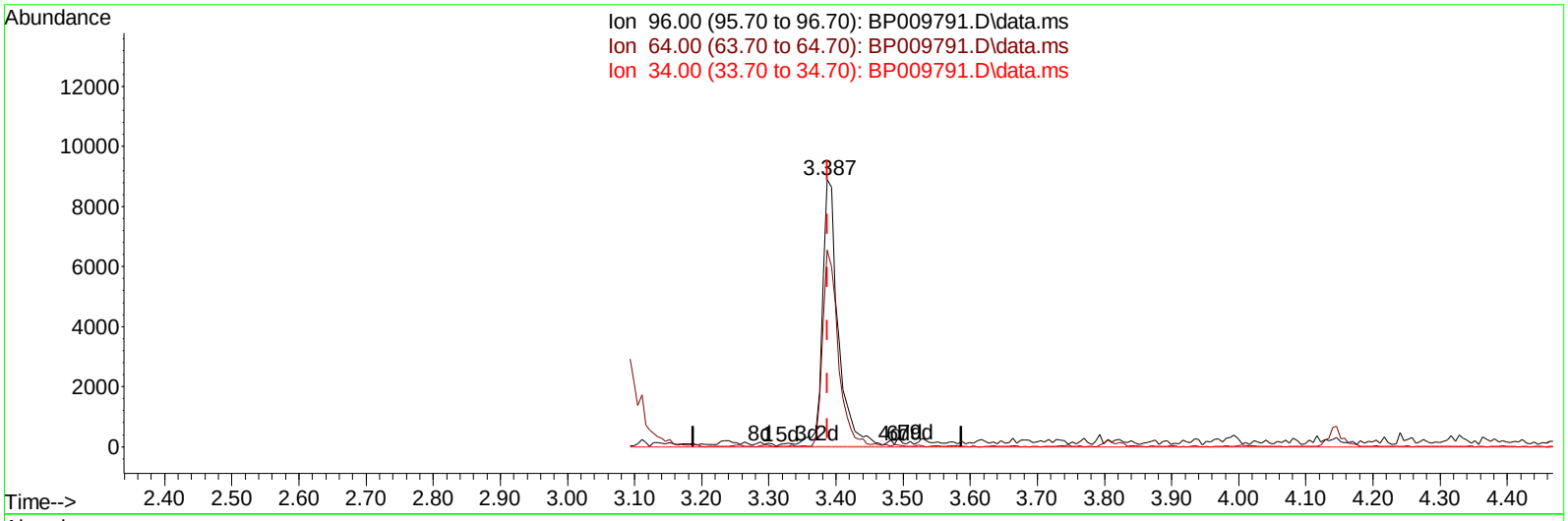

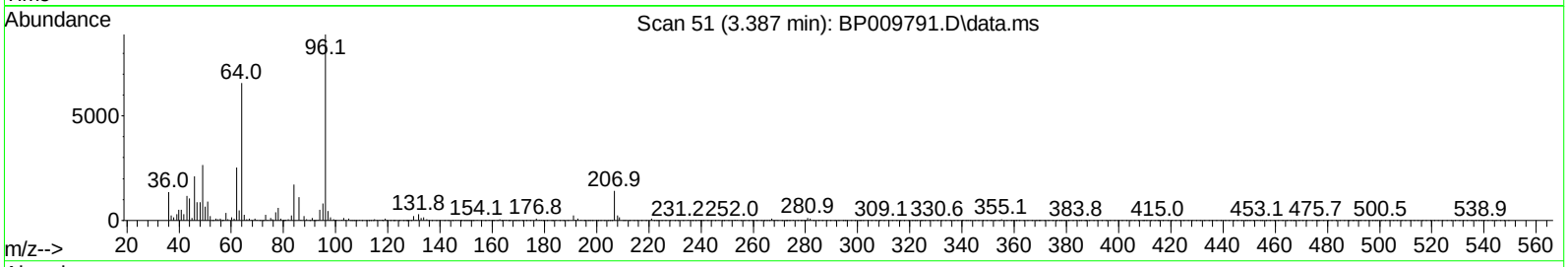

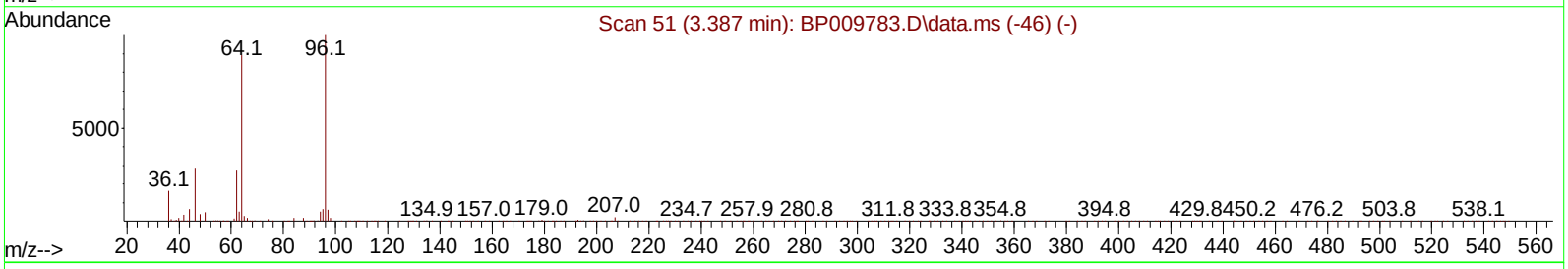

TIC: BP009791.D\data.ms

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

3.387min (+ 0.000) 5.17 ng/uL m

response 14312

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.80#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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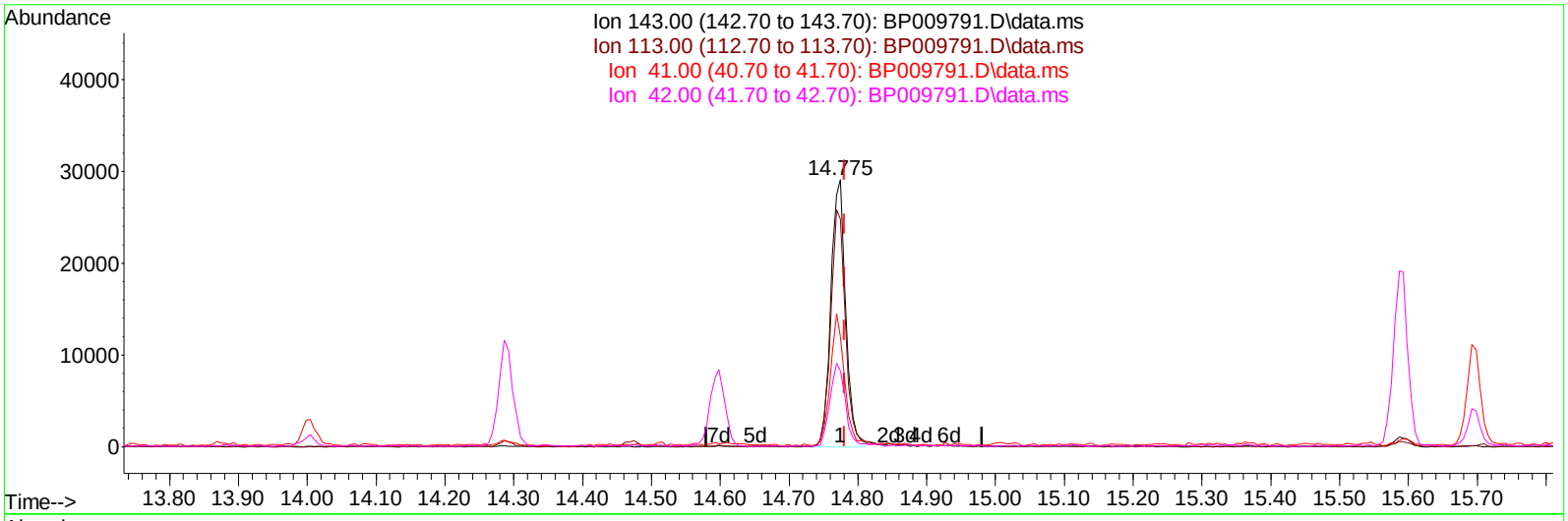

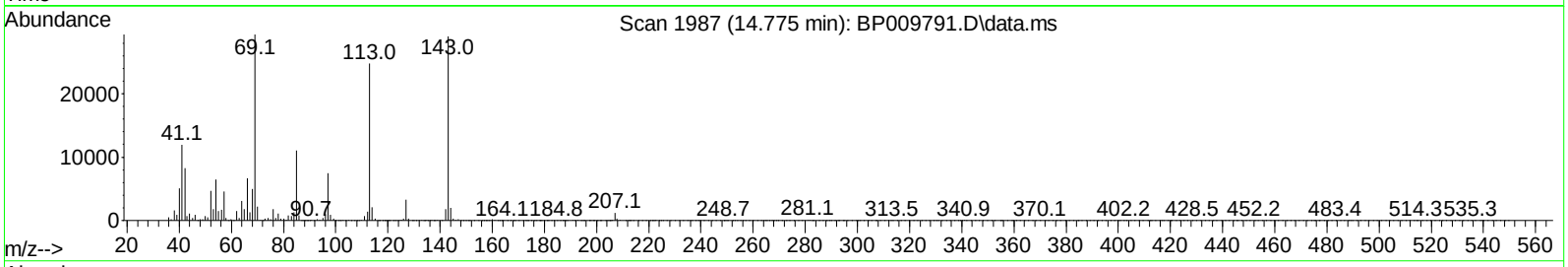

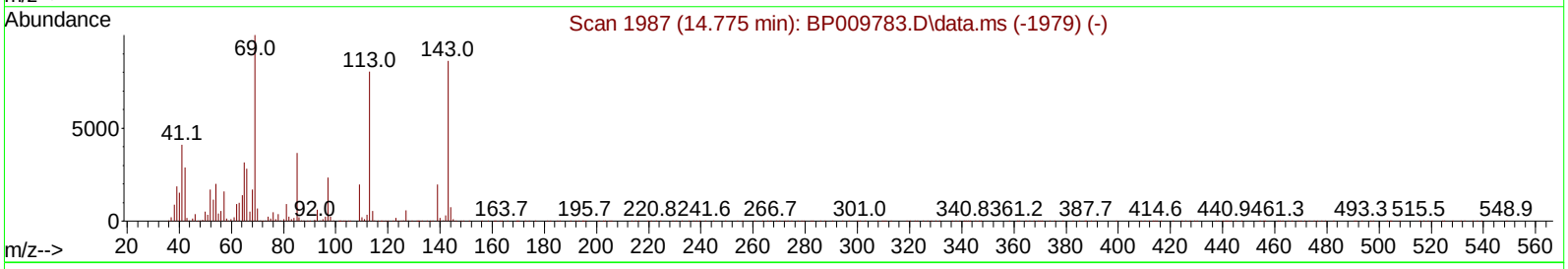

TIC: BP009791.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 8.24 ng/ul

response 42917

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.17#
41.00	23.20	41.19#
42.00	18.00	28.52#

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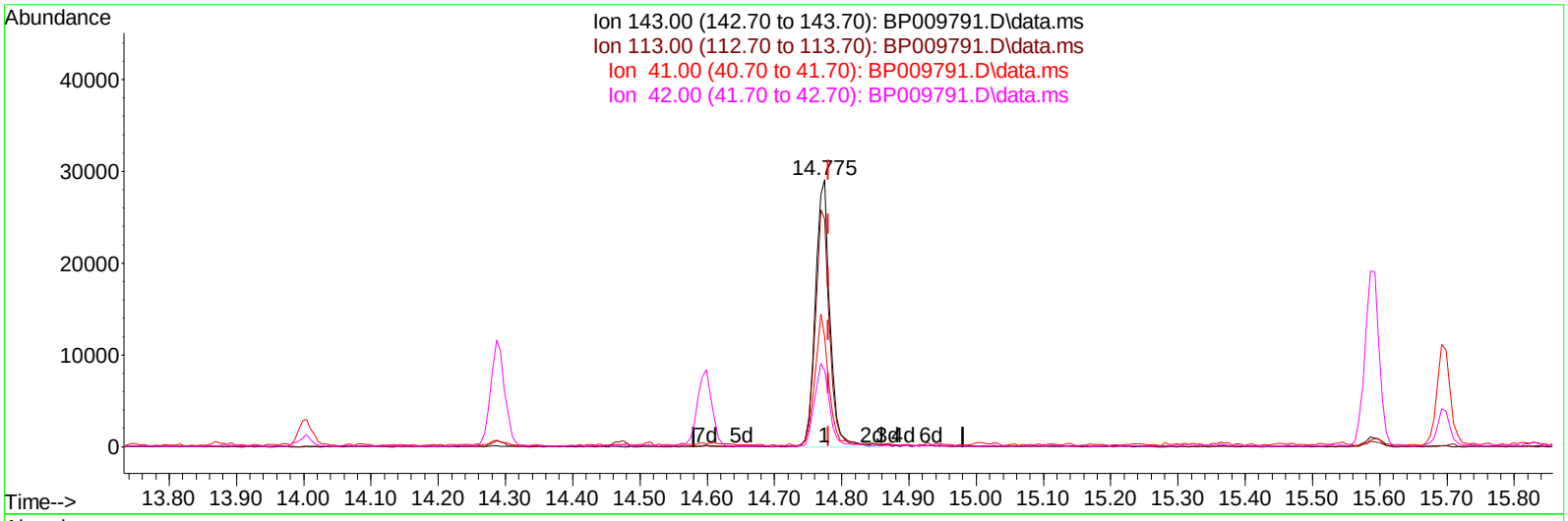

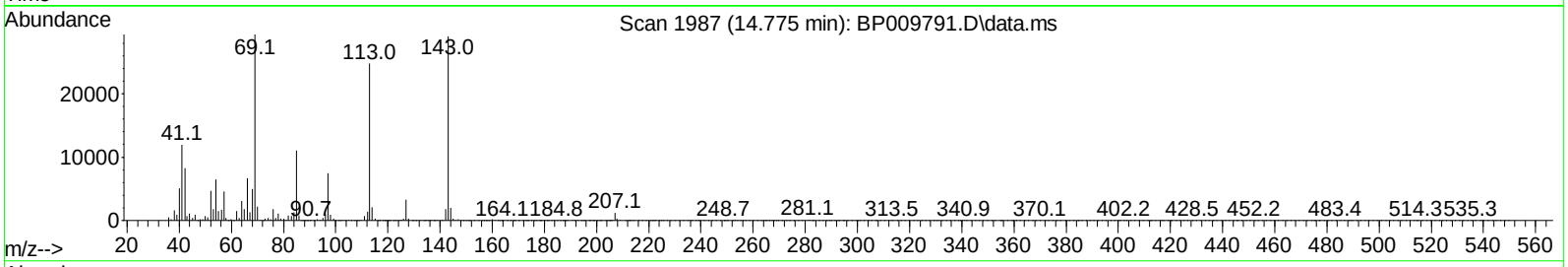

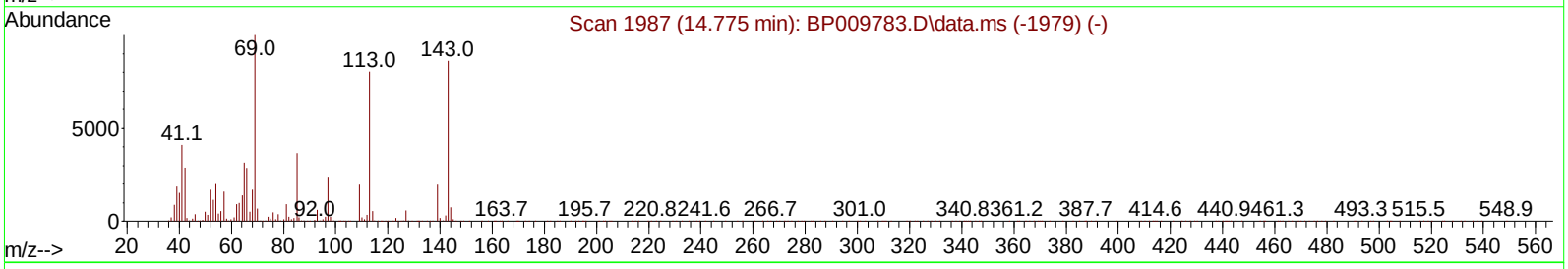

TIC: BP009791.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 8.39 ng/ul m

response 43721

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.17#
41.00	23.20	41.19#
42.00	18.00	28.52#

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	8.016	152	123470	20.000	ng/ul		0.05
20) Naphthalene-d8	10.781	136	509129	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.598	164	365676	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.345	188	838591	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.427	240	911133	20.000	ng/ul	#	0.00
88) Perylene-d12	23.892	264	881180	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.387	96	14312m	5.173	ng/uL		0.00
4) Pyridine-d5	3.811	84	89211	11.696	ng/ul		0.00
7) Phenol-d5	7.128	99	91357	8.428	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	392120	60.762	ng/ul		0.02
11) 2-Chlorophenol-d4	7.505	132	238573	29.177	ng/ul		0.00
15) 4-Methylphenol-d8	8.669	113	169596	18.742	ng/ul		0.00
21) Nitrobenzene-d5	9.128	128	155292	42.297	ng/ul		0.00
24) 2-Nitrophenol-d4	9.851	143	163028	40.502	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.393	165	286170	33.975	ng/ul		0.00
31) 4-Chloroaniline-d4	10.904	131	266526	22.216	ng/ul		0.00
46) Dimethylphthalate-d6	14.004	166	1128988	39.304	ng/ul		0.00
49) Acenaphthylene-d8	14.292	160	1274625	37.456	ng/ul		0.00
54) 4-Nitrophenol-d4	14.775	143	43721m	8.393	ng/ul		0.00
60) Fluorene-d10	15.592	176	991960	41.051	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	165359	33.286	ng/ul		0.00
73) Anthracene-d10	17.445	188	1641120	40.874	ng/ul		0.00
81) Pyrene-d10	19.674	212	2062962	41.457	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.733	264	1961491	42.745	ng/ul		0.00
Target Compounds							Qvalue
8) Phenol	7.158	94	33409	3.079	ng/ul	#	92
12) 2-Chlorophenol	7.540	128	21886	2.655	ng/ul		96
32) 4-Chloroaniline	10.928	127	39617	3.337	ng/ul		96

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

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