Quantitation Report (Qedit)

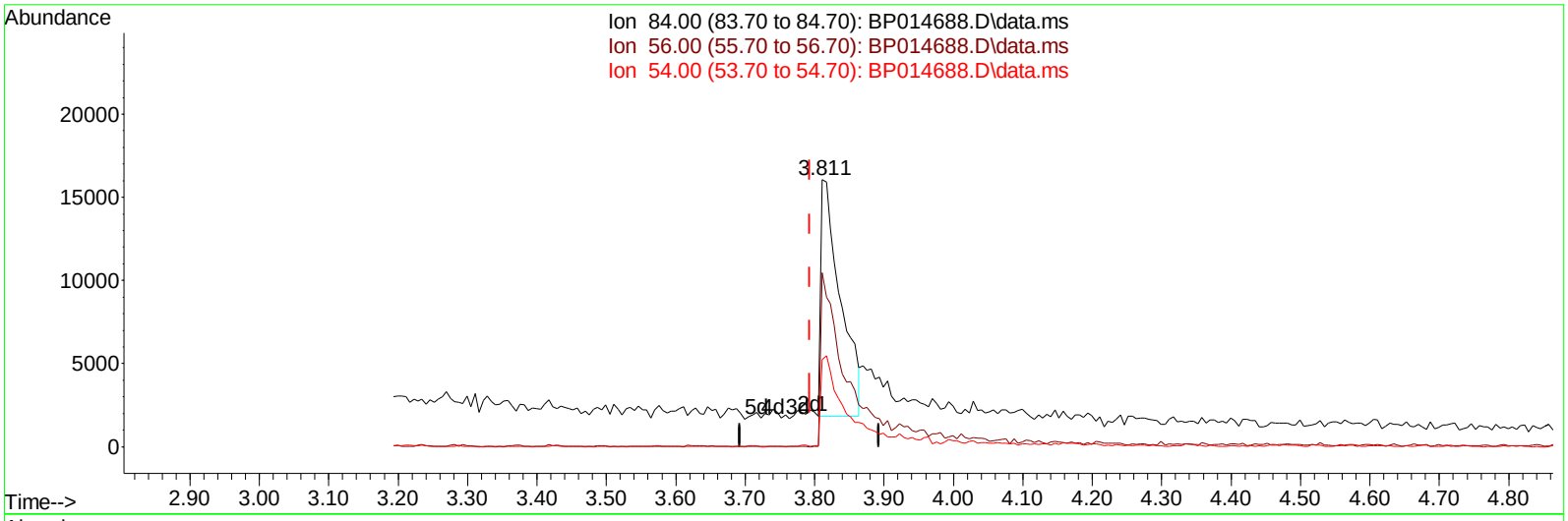
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041223\
Data File : BP014688.D
Acq On : 13 Apr 2023 00:07
Operator : CG/JU
Sample : 02250-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

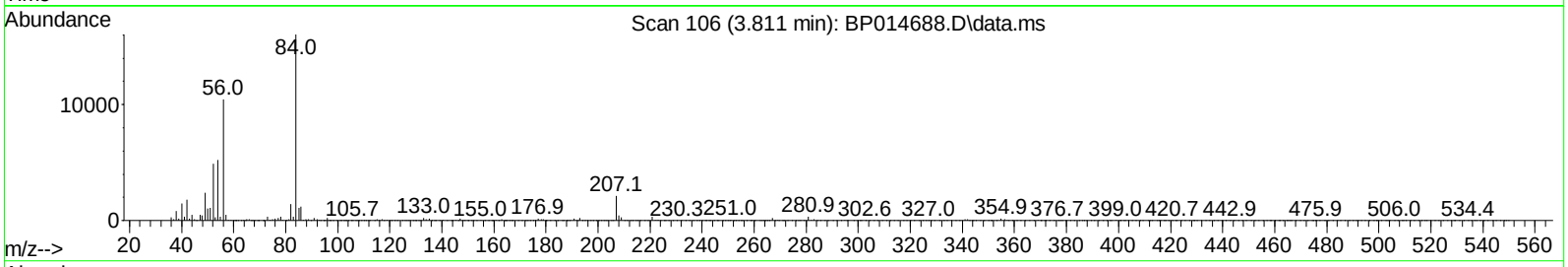
Quant Time: Apr 13 02:06:08 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 13 02:01:50 2023
Response via : Initial Calibration

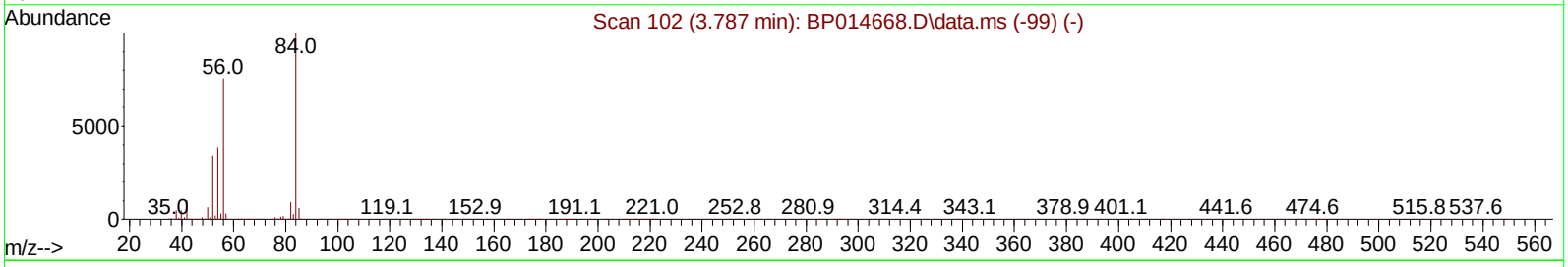
Instrument :
BNA_P
ClientSampleId :
BHDY3

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/13/2023
Supervised By :Jagrut Upadhyay 04/13/2023







TIC: BP014688.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.018) 4.70 ng/ul

response 28177

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.17
54.00	32.60	32.58
0.00	0.00	0.00

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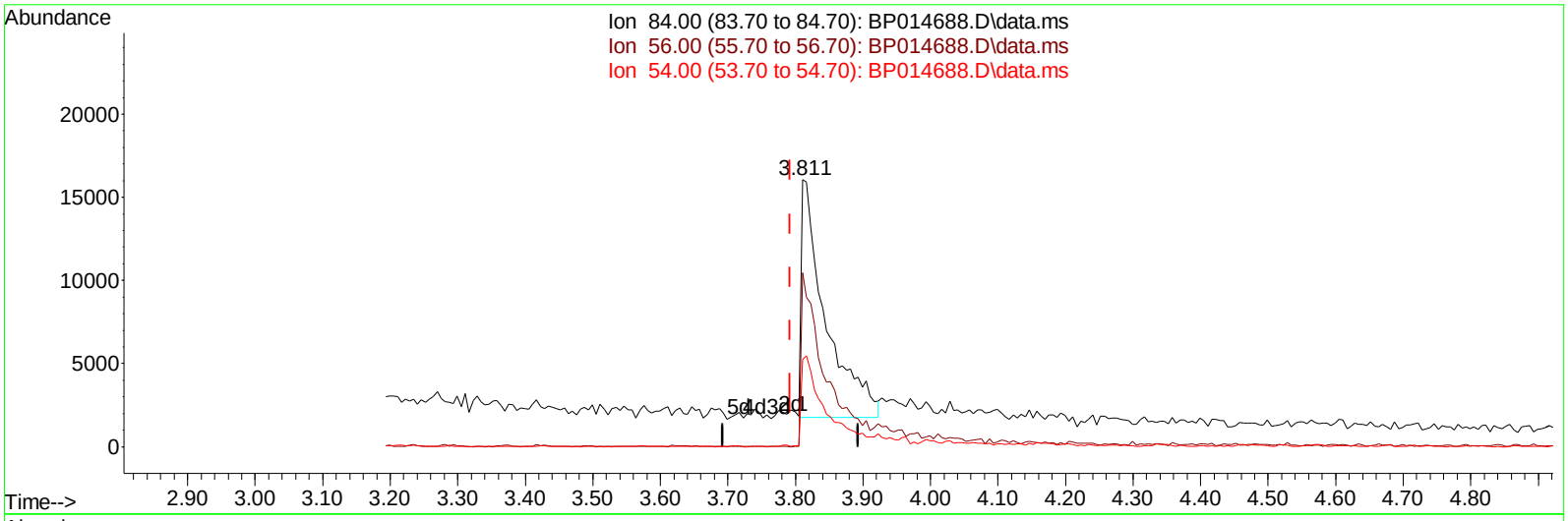

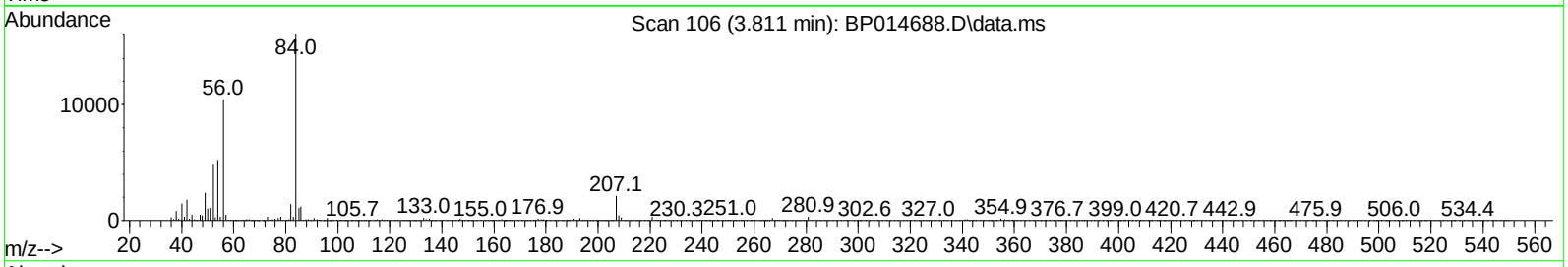

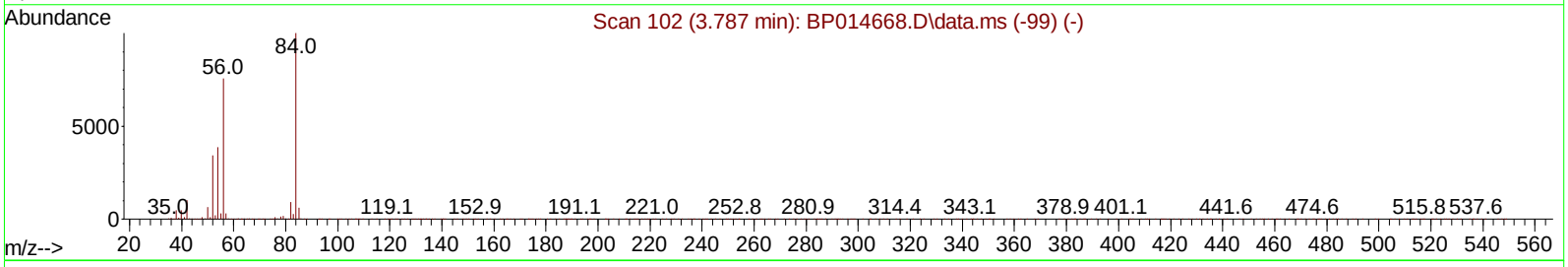

TIC: BP014688.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.018) 5.97 ng/ul m

response 35756

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.17
54.00	32.60	32.58
0.00	0.00	0.00

Quantitation Report (Qedit)

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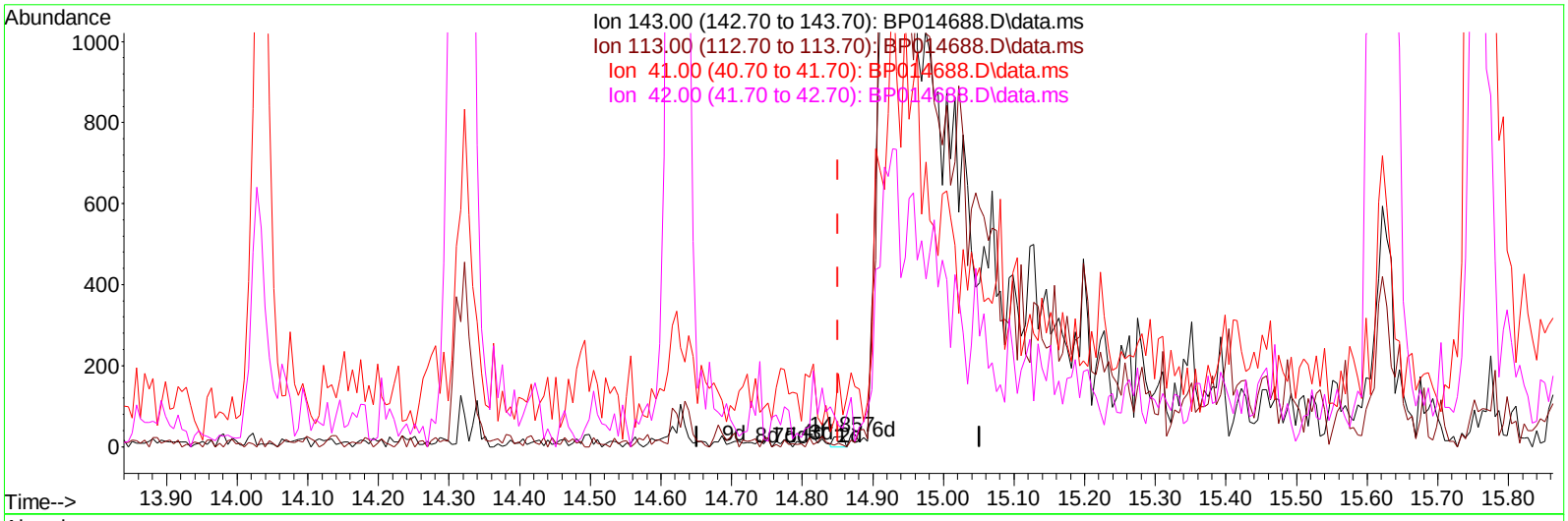

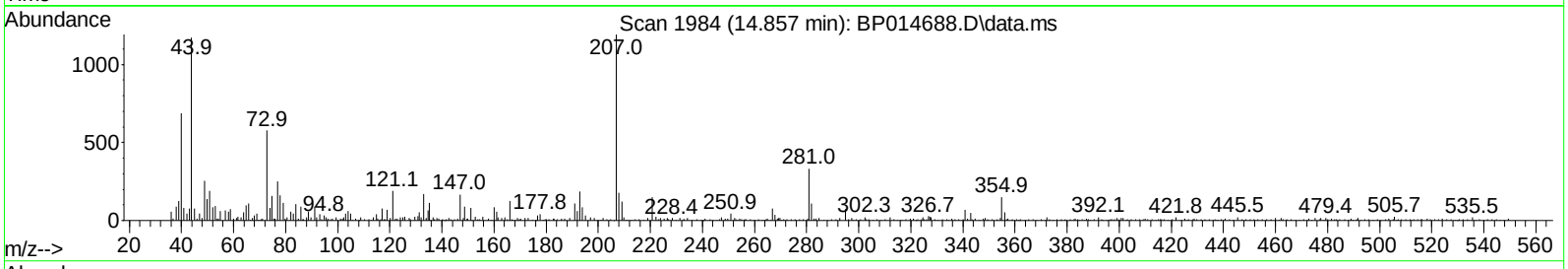

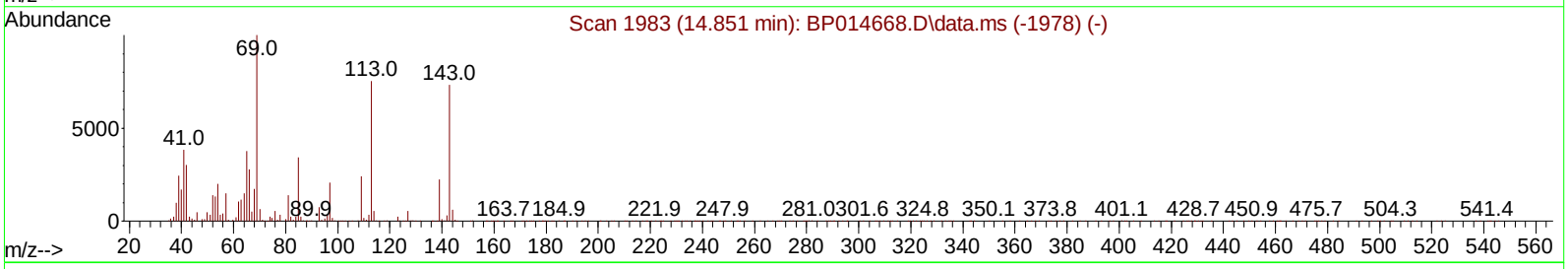

TIC: BP014688.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (+ 0.006) 0.01 ng/ul

response 26

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.00
41.00	44.20	316.00#
42.00	31.50	184.00#

Quantitation Report (Qedit)

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ClientSampleId :
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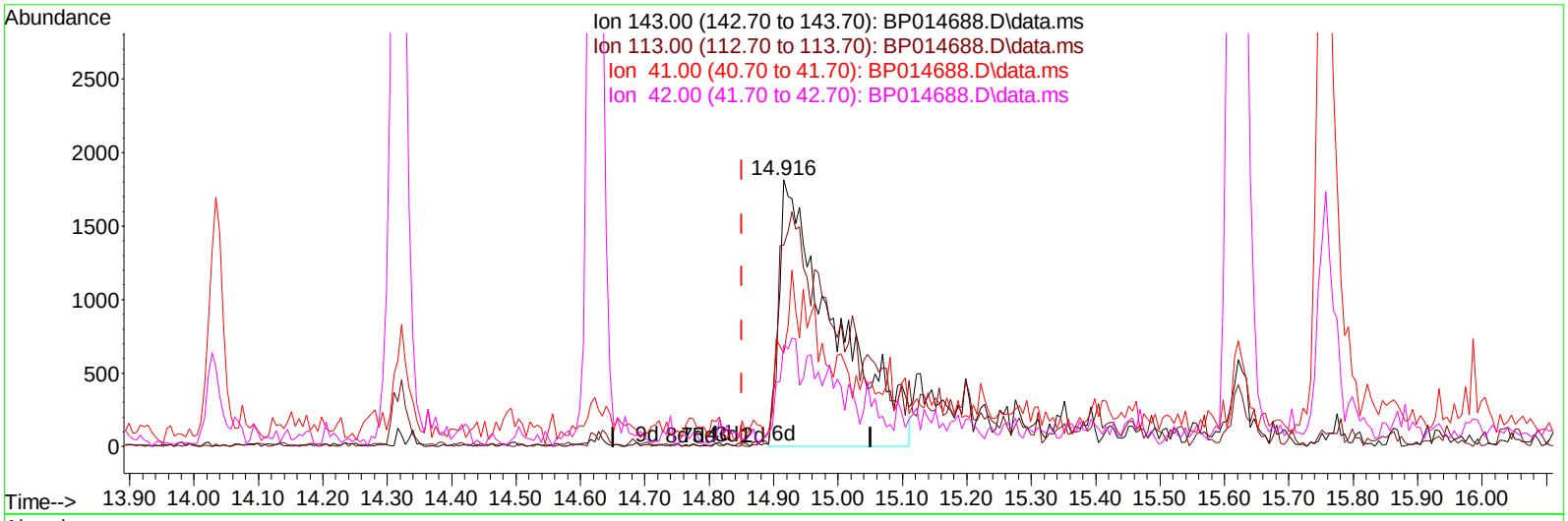

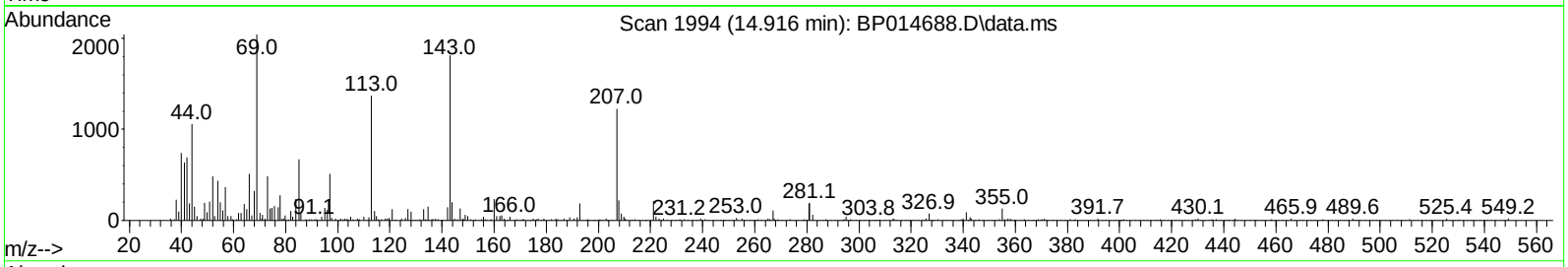

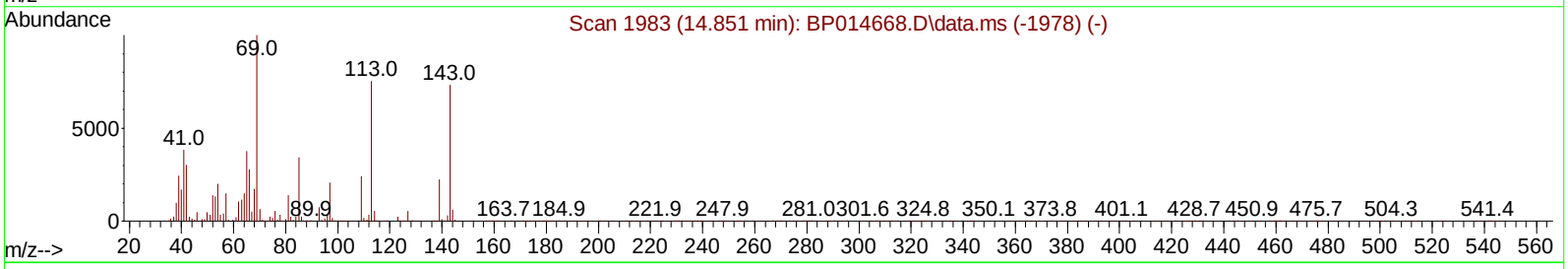

TIC: BP014688.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.065) 4.50 ng/ul m

response 10695

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	75.65#
41.00	44.20	34.99#
42.00	31.50	38.07#

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	7.969	152	90688	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	353126	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.622	164	193965	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.387	188	393711	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	382417	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	460009	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	11587	4.955	ng/uL	0.00
4) Pyridine-d5	3.811	84	35756m	5.966	ng/ul	0.02
7) Phenol-d5	7.152	99	52936	6.329	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.299	67	176210	32.837	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	149591	24.746	ng/ul	0.00
15) 4-Methylphenol-d8	8.693	113	94695	14.036	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	87787	30.795	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	96599	29.593	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	147486	24.809	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	171951	21.959	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	533787	33.941	ng/ul	0.00
49) Acenaphthylene-d8	14.316	160	580553	33.268	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	10695m	4.501	ng/ul	0.07
60) Fluorene-d10	15.622	176	457944	34.917	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	77513	26.756	ng/ul	0.00
73) Anthracene-d10	17.487	188	719545	37.764	ng/ul	0.00
81) Pyrene-d10	19.722	212	855399	33.301	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	934988	38.256	ng/ul	0.00

Target Compounds Qvalue

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

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