Quantitation Report (Qedit)

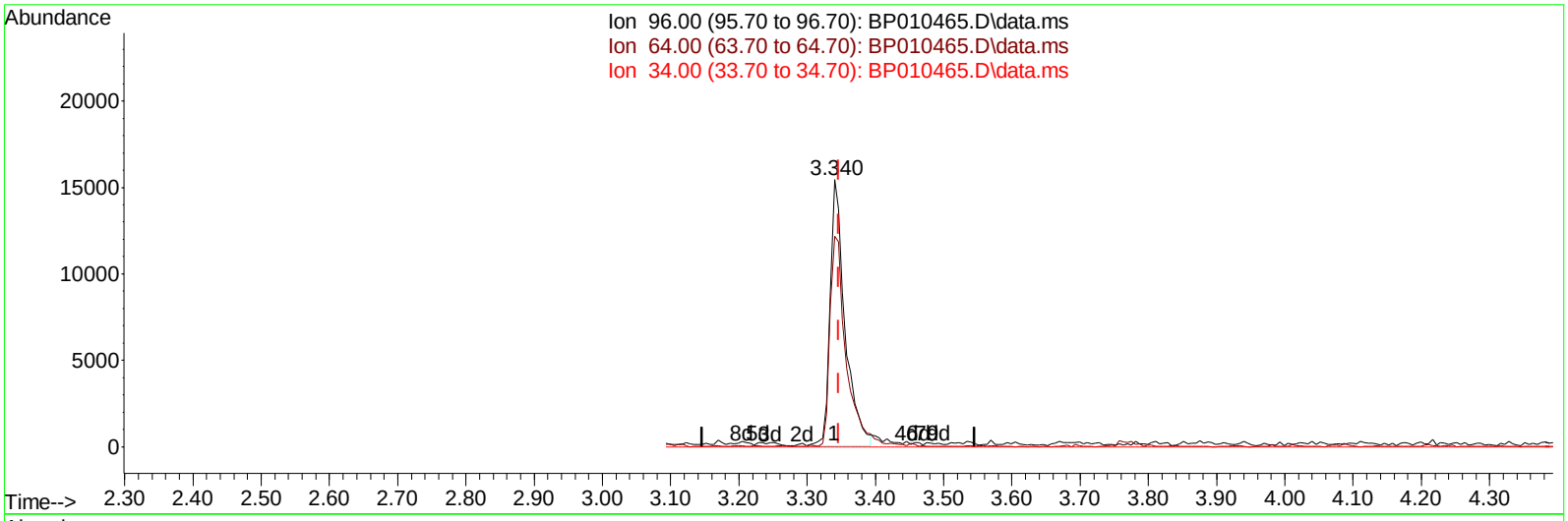
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
Data File : BP010465.D
Acq On : 21 May 2022 08:41
Operator : CG/JU
Sample : N2918-17
Misc :
ALS Vial : 40 Sample Multiplier: 1

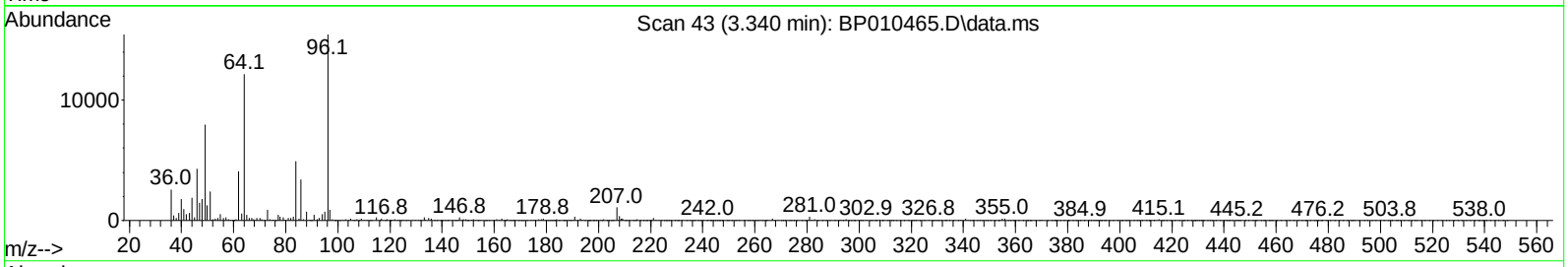
Quant Time: May 23 00:43:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

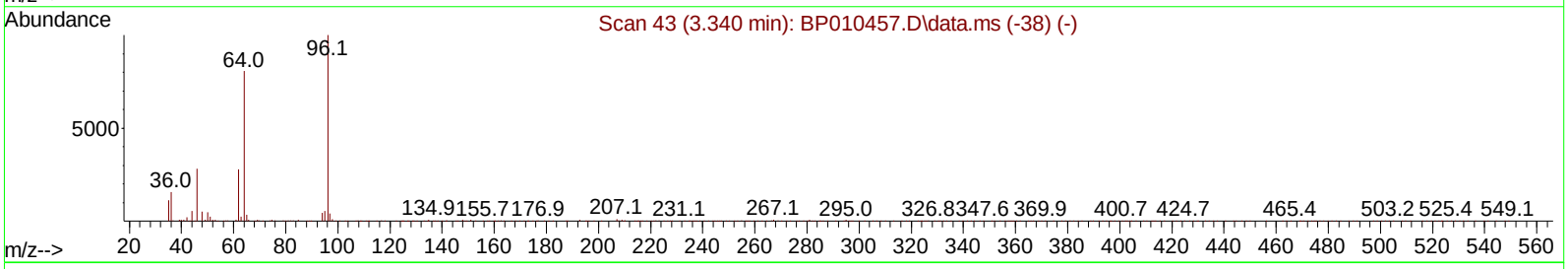
Instrument :
BNA_P
ClientSampleId :
DBTD3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/23/2022
Supervised By :mohammad ahmed 05/25/2022







TIC: BP010465.D\data.ms

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

3.340min (-0.006) 4.65 ng/uL

response 23269

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.80#
34.00	0.00	0.00
0.00	0.00	0.00

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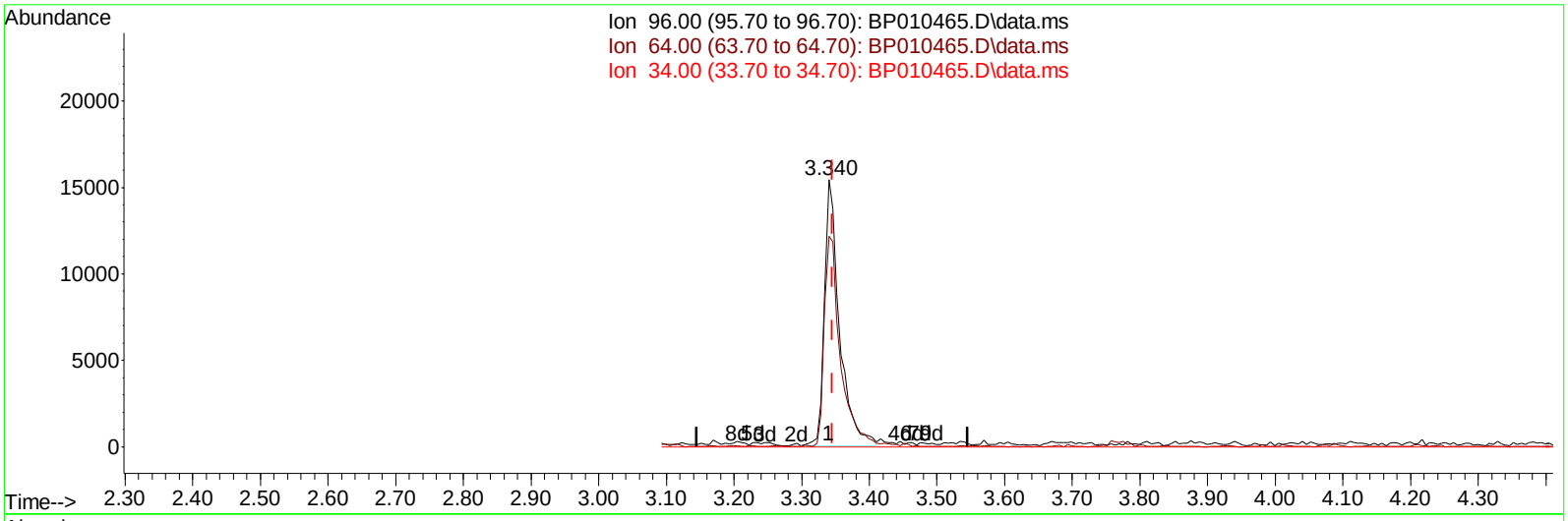

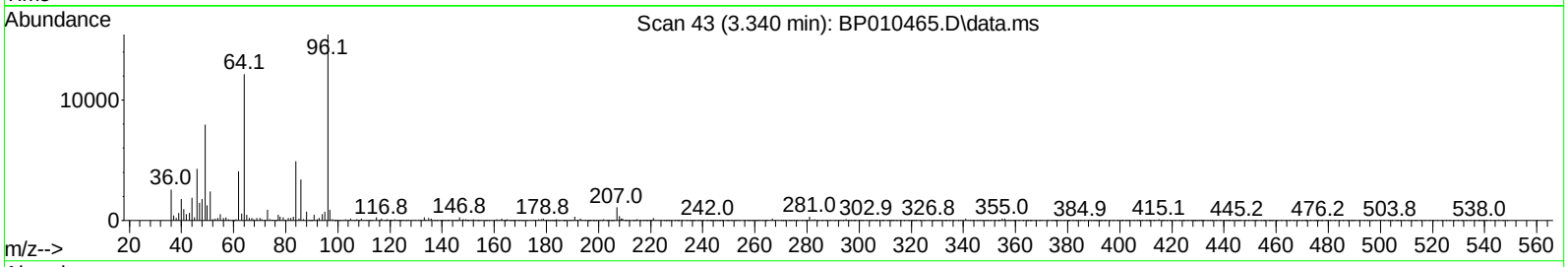

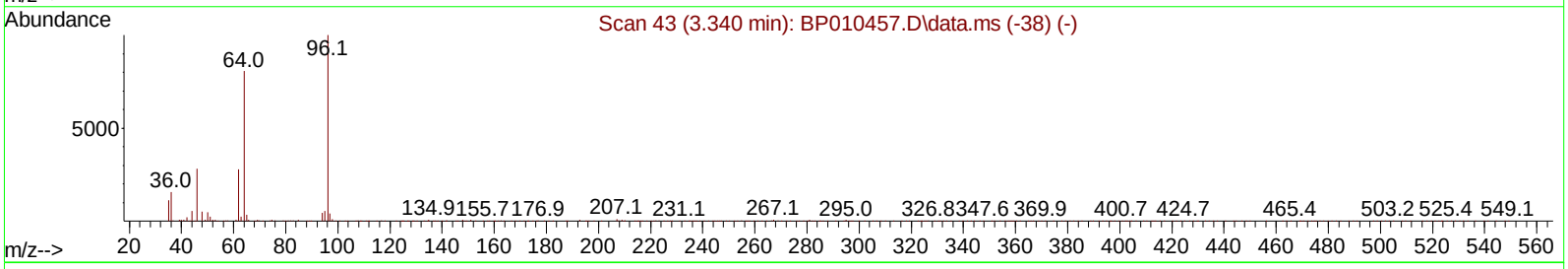

TIC: BP010465.D\data.ms

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

3.340min (-0.006) 4.77 ng/uL m

response 23846

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.80#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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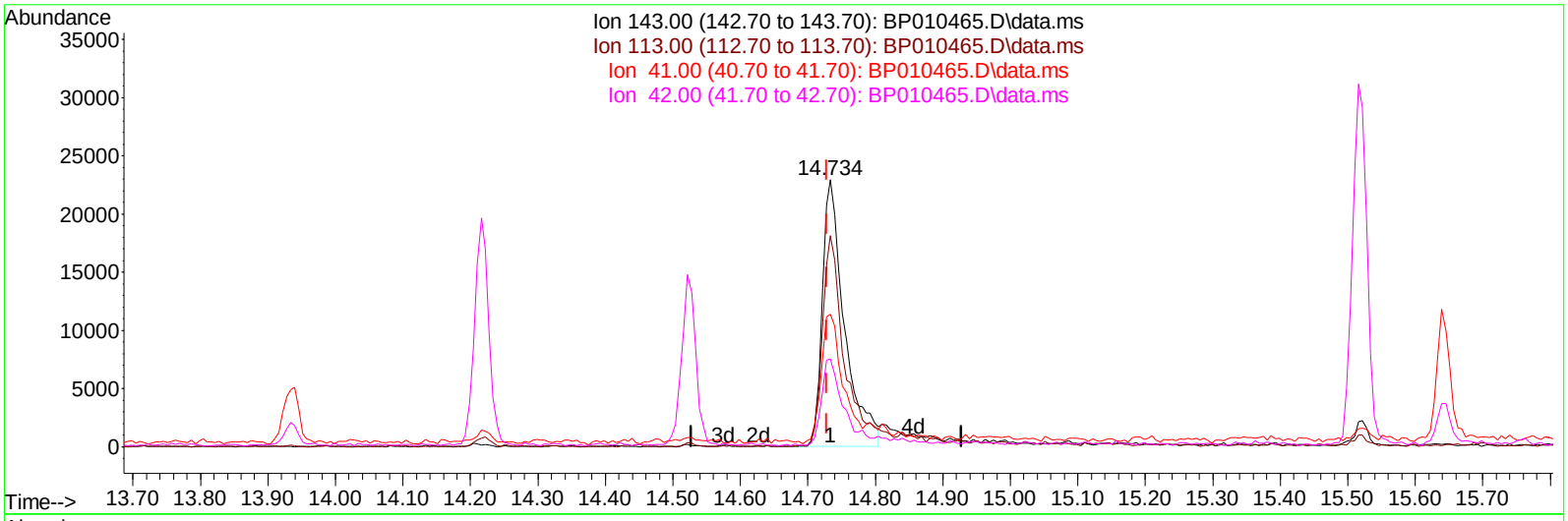

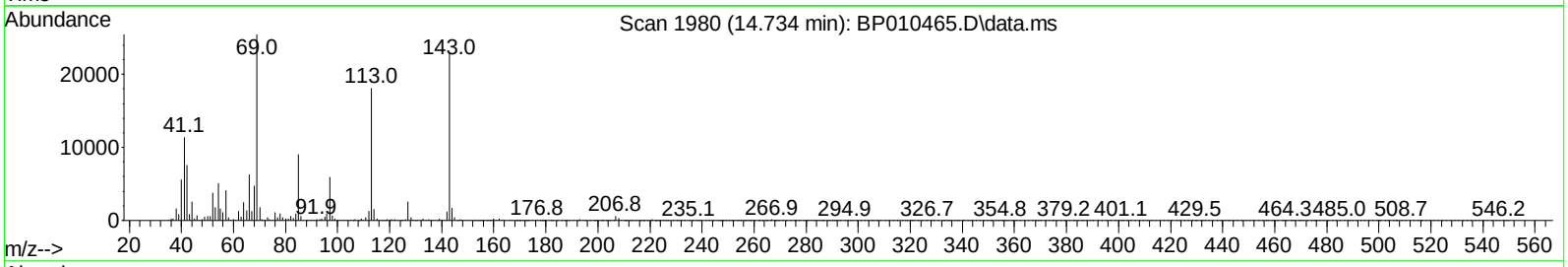

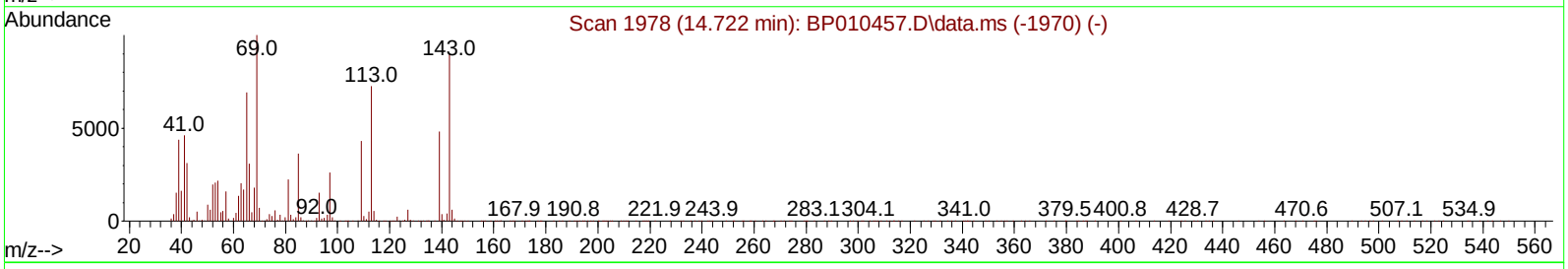

TIC: BP010465.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 6.37 ng/ul

response 52632

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.11#
41.00	23.20	49.76#
42.00	18.00	32.98#

Quantitation Report (Qedit)

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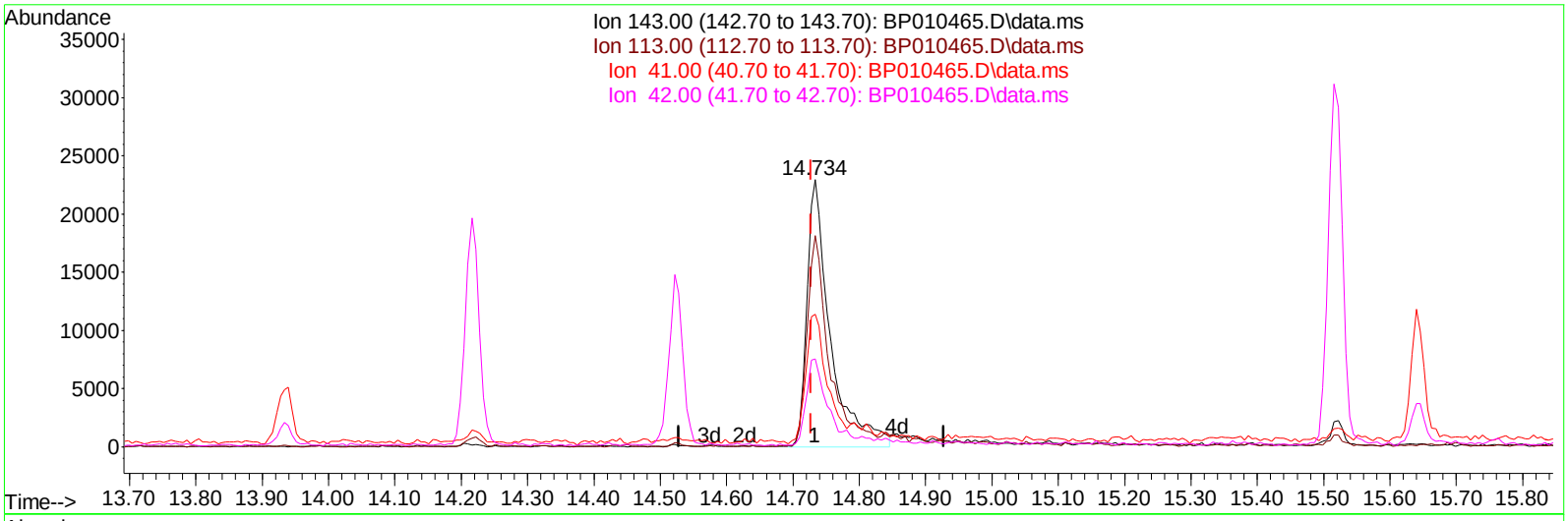

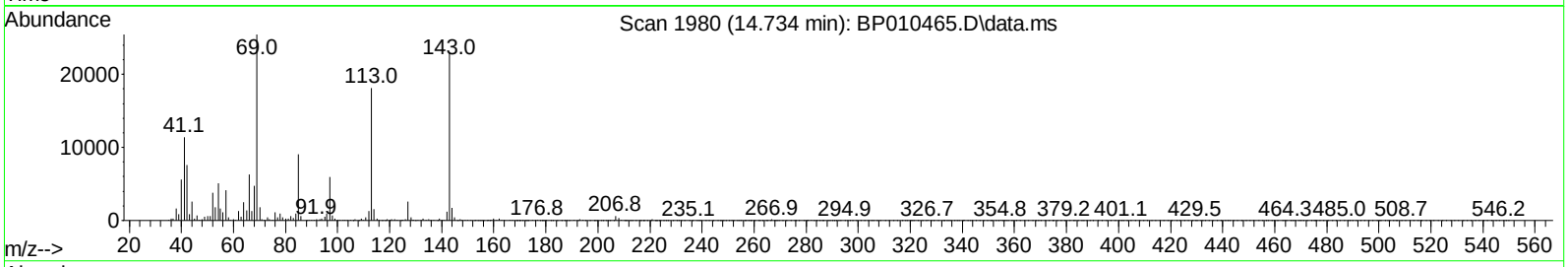

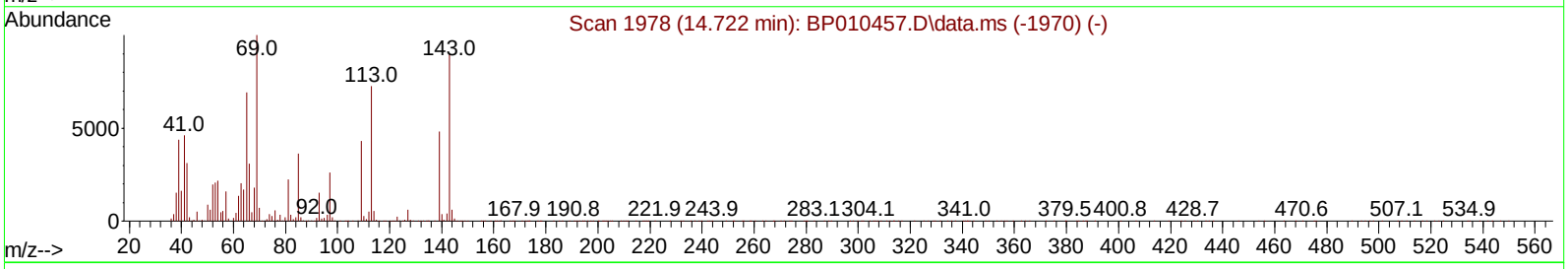

TIC: BP010465.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 6.81 ng/ul m

response 56234

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.11#
41.00	23.20	49.76#
42.00	18.00	32.98#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	213268	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	949158	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.522	164	629975	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1290270	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	932325	20.000	ng/ul	# 0.00
88) Perylene-d12	23.786	264	760245	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	23846m	4.766	ng/uL	0.00
4) Pyridine-d5	3.764	84	139541	9.822	ng/ul	0.00
7) Phenol-d5	7.057	99	151883	8.485	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	461912	38.408	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	441910	32.263	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	310338	20.401	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	302419	41.169	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	323283	41.527	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	523554	35.567	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	632510	28.899	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1917491	41.312	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	2214003	41.354	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	56234m	6.807	ng/ul	0.00
60) Fluorene-d10	15.522	176	1649584	40.852	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	159806	21.178	ng/ul	0.00
73) Anthracene-d10	17.380	188	2648201	45.326	ng/ul	0.00
81) Pyrene-d10	19.610	212	2866261	57.889	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1756858	46.022	ng/ul	0.00

Target Compounds Qvalue

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

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