Quantitation Report (Qedit)

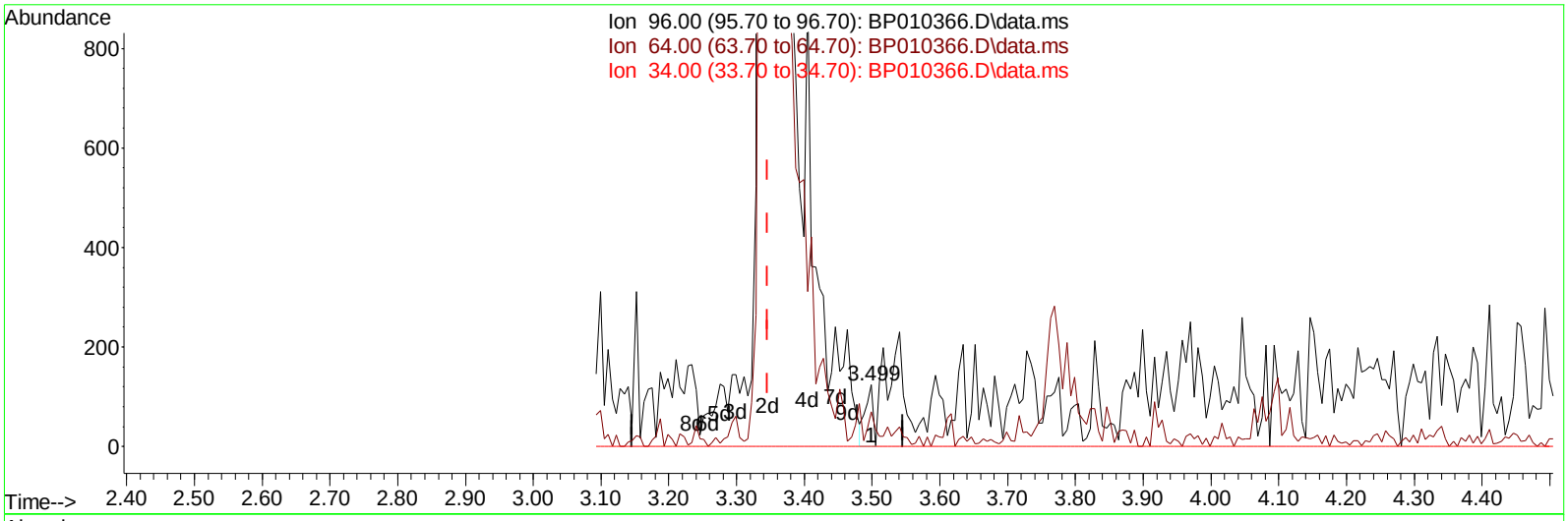
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010366.D
Acq On : 16 May 2022 20:20
Operator : CG/JU
Sample : N2746-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

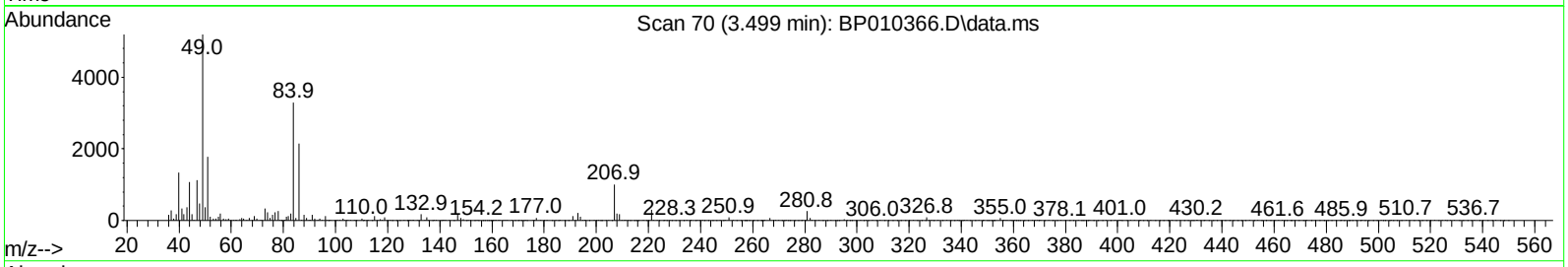
Quant Time: May 18 01:20:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

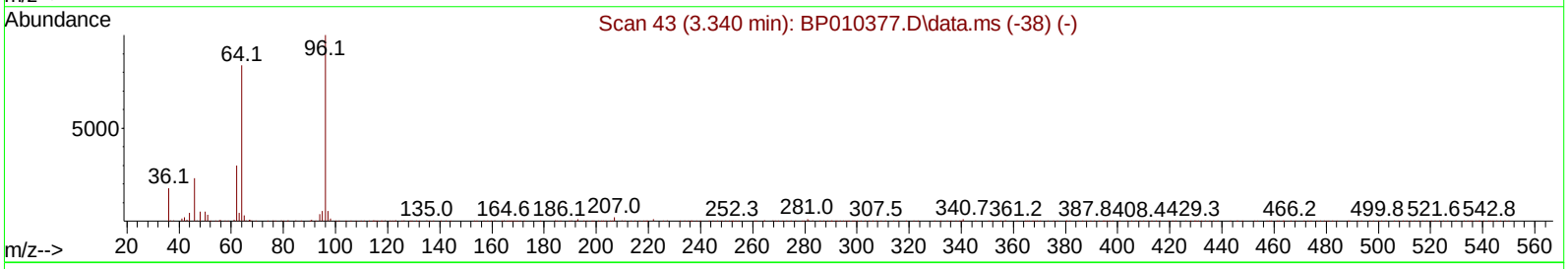
Instrument :
BNA_P
ClientSampleId :
DBSL1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/18/2022
Supervised By :mohammad ahmed 05/18/2022







TIC: BP010366.D\data.ms

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

3.499min (+ 0.153) 0.05 ng/uL

response 97

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.45
34.00	0.00	0.00
0.00	0.00	0.00

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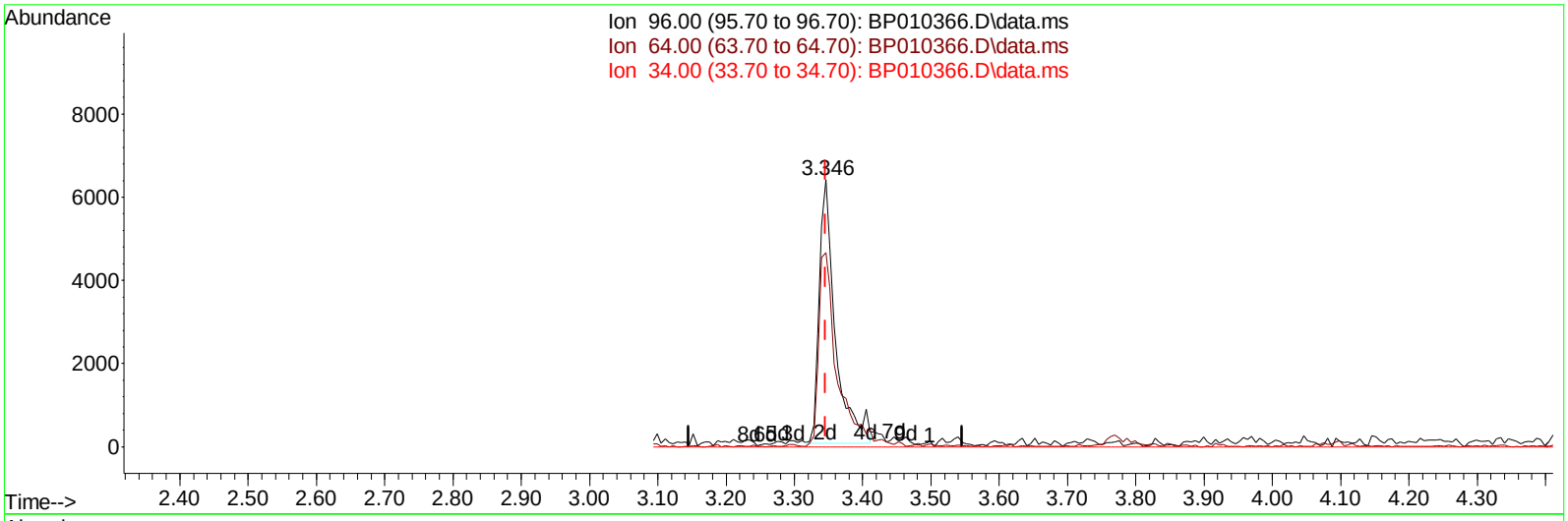

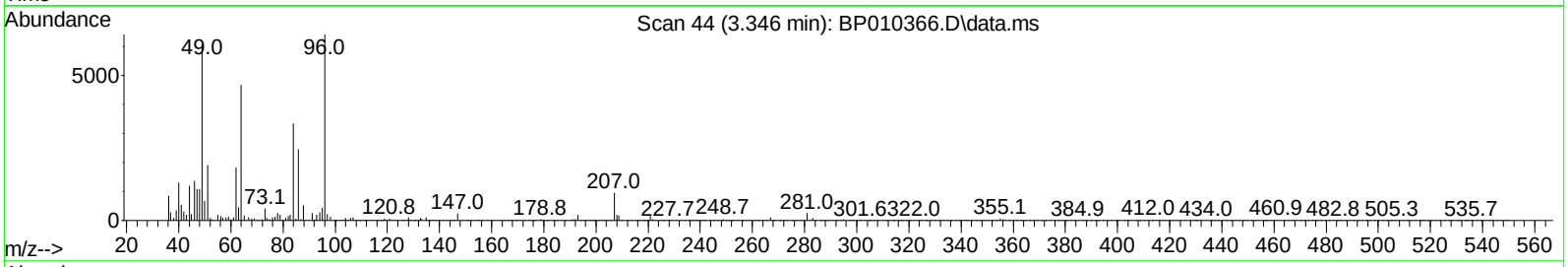

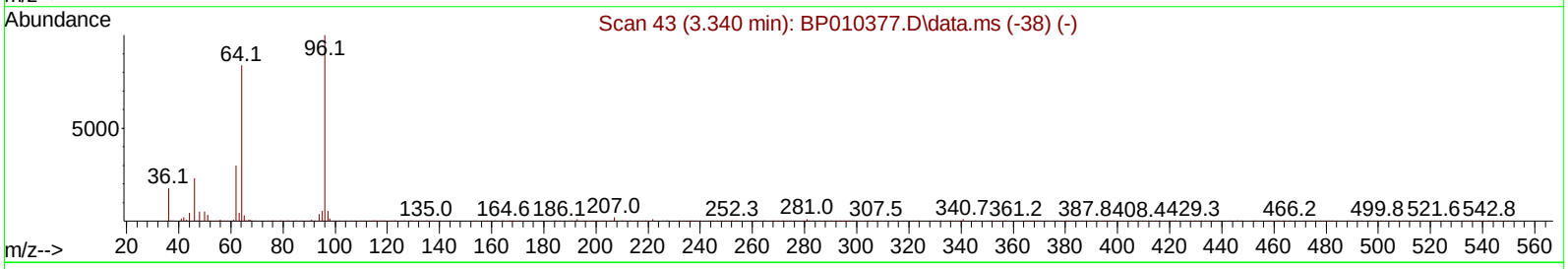

TIC: BP010366.D\data.ms

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

3.346min (-0.000) 5.08 ng/uL m

response 10433

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	72.85#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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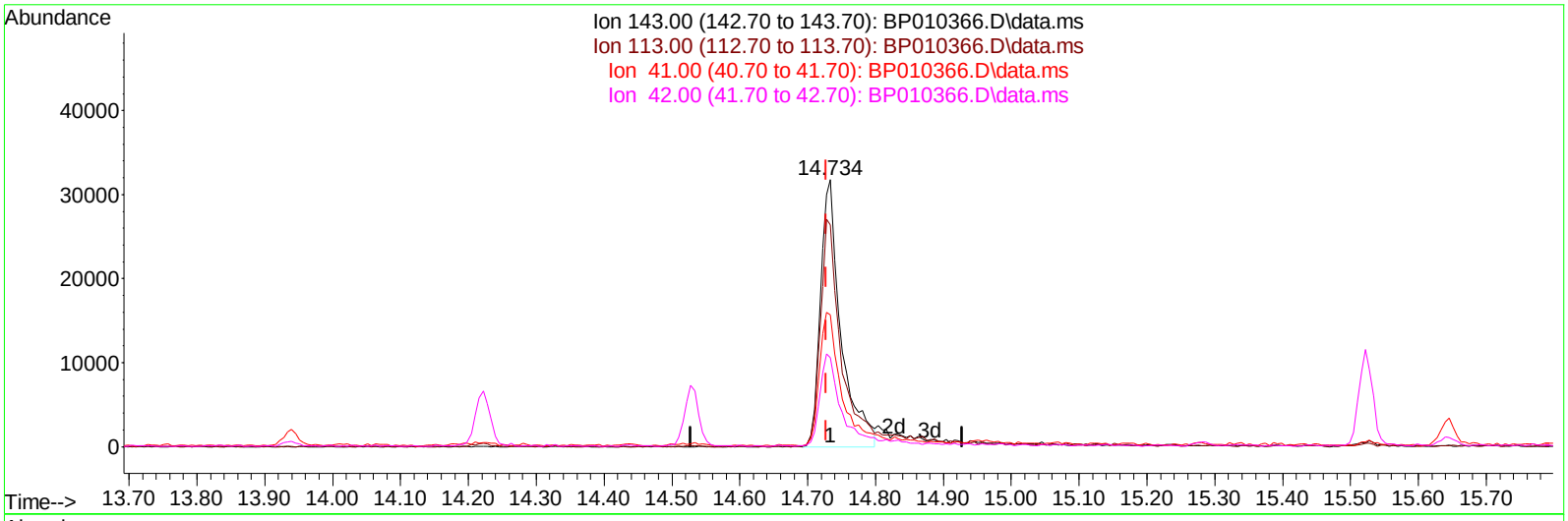

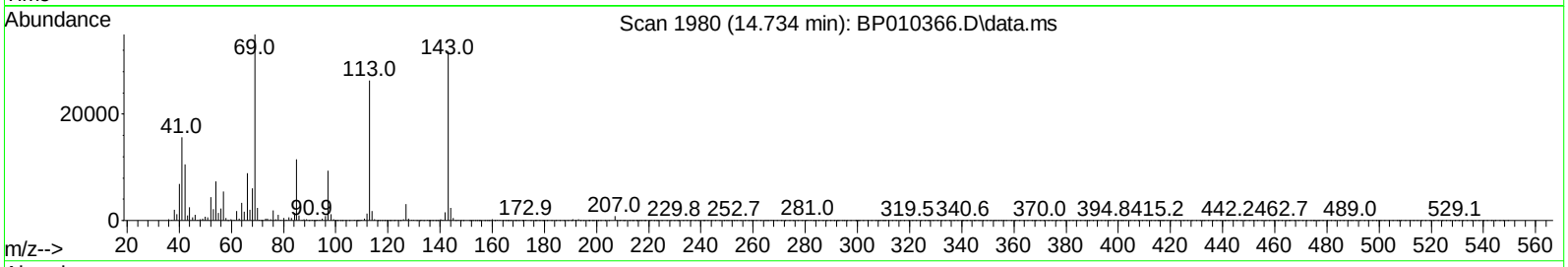

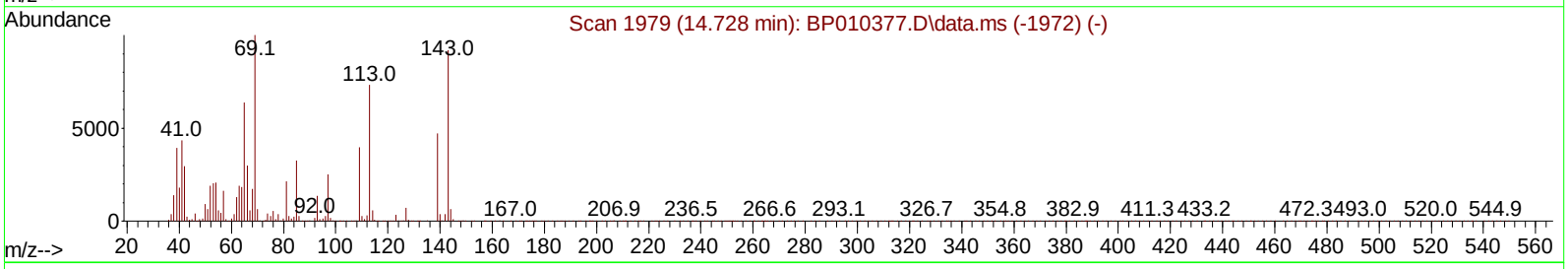

TIC: BP010366.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 18.44 ng/ul

response 66615

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.07#
41.00	23.20	49.30#
42.00	18.00	33.26#

Quantitation Report (Qedit)

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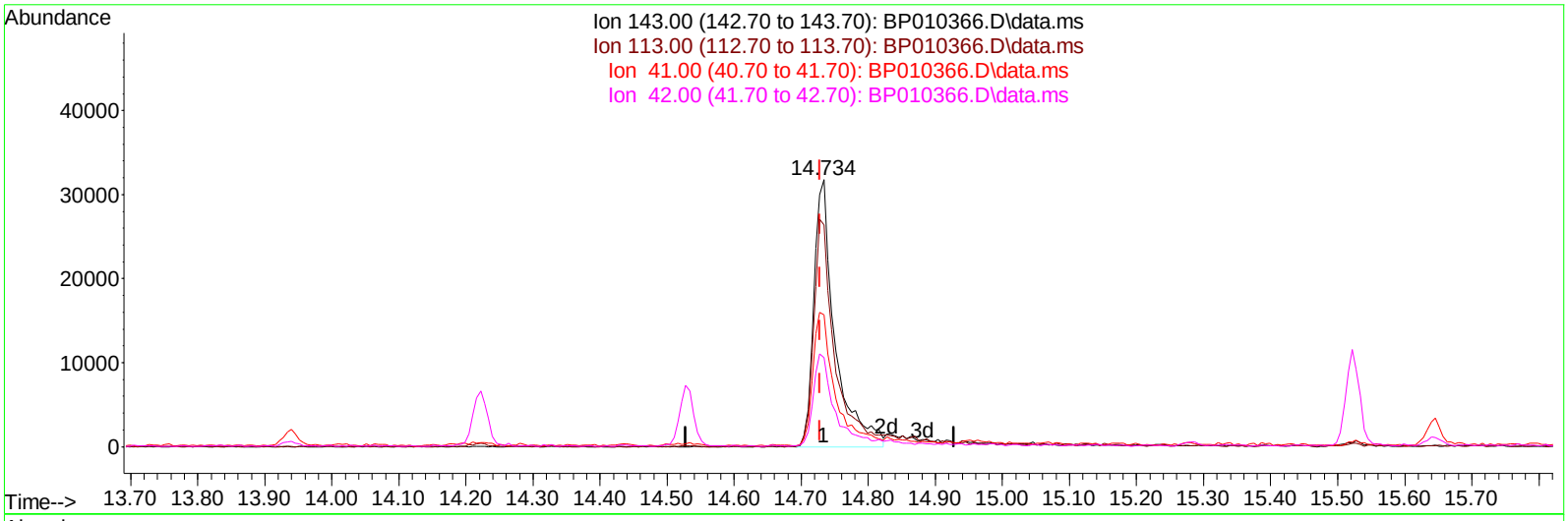

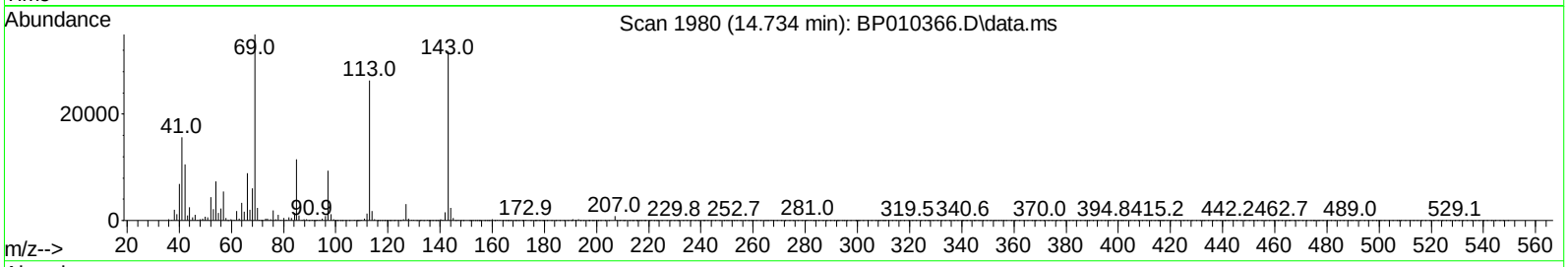

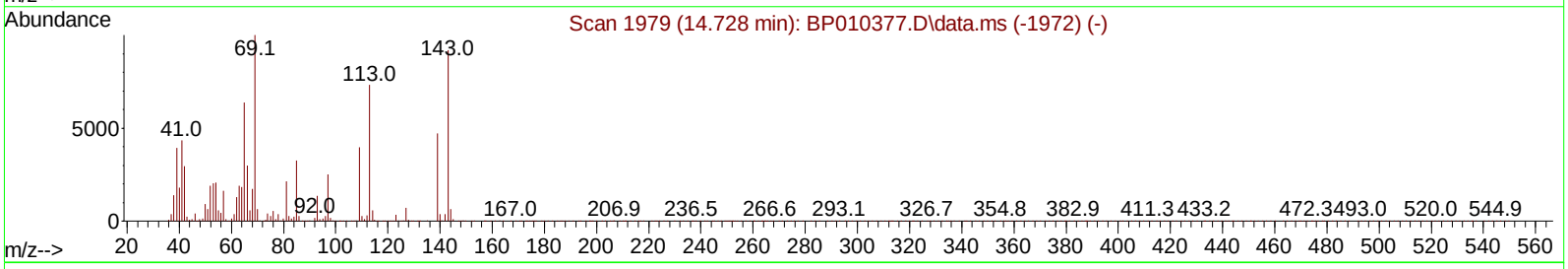

TIC: BP010366.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 19.19 ng/ul m

response 69320

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.07#
41.00	23.20	49.30#
42.00	18.00	33.26#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	87625	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	389153	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	275472	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	625364	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	682627	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	745785	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	10433m	5.075	ng/uL	0.00
4) Pyridine-d5	3.764	84	135369	23.191	ng/ul	0.00
7) Phenol-d5	7.063	99	194338	26.424	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	137199	27.766	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	156166	27.750	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	173144	27.702	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	86504	28.722	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	86155	26.993	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	159017	26.348	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	244893	27.291	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	588600	29.001	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	672794	28.739	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	69320m	19.189	ng/ul	0.00
60) Fluorene-d10	15.522	176	521117	29.513	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	43511	11.897	ng/ul	0.00
73) Anthracene-d10	17.381	188	852898	30.119	ng/ul	0.00
81) Pyrene-d10	19.616	212	1084884	29.926	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1133596	30.271	ng/ul	0.00
Target Compounds						Qvalue

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

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