

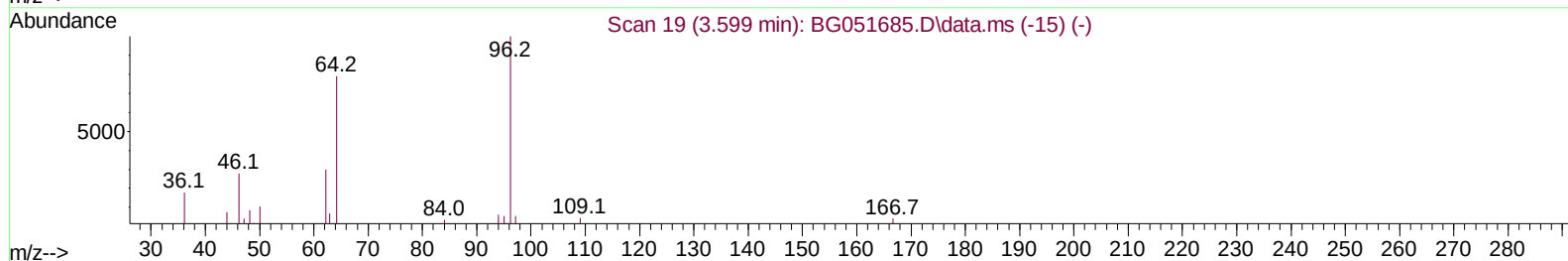
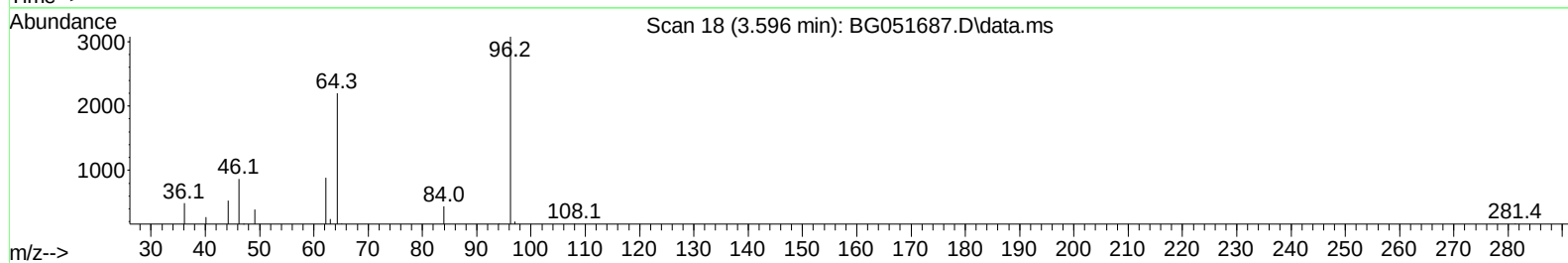
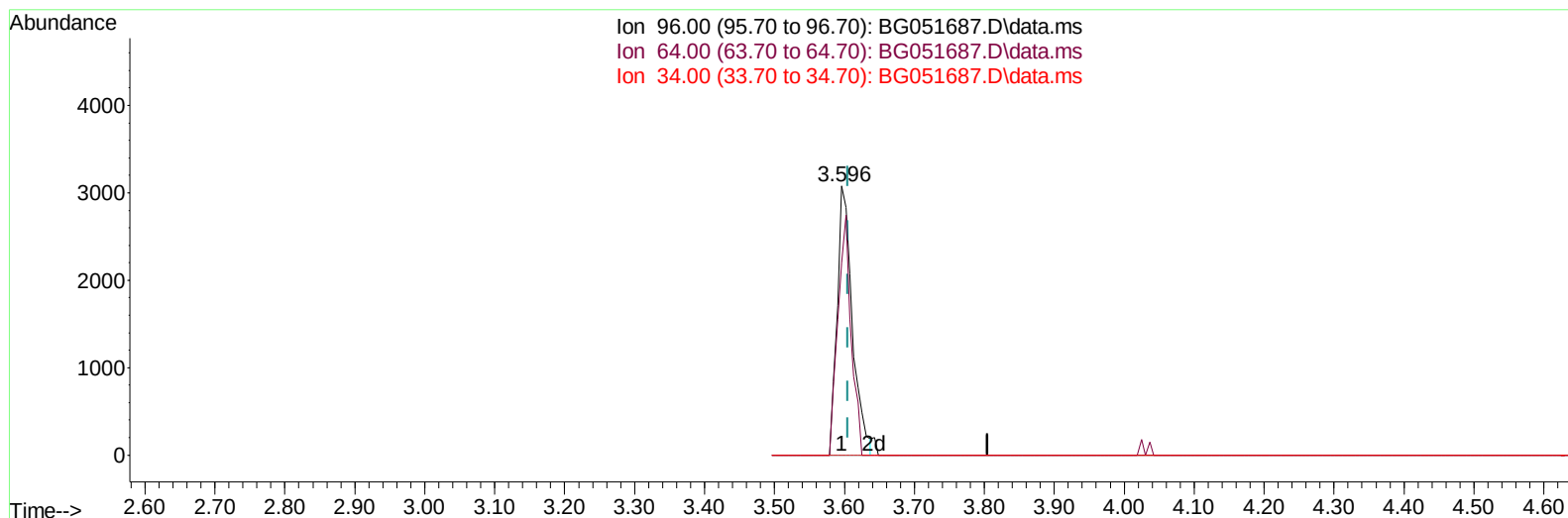
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
Data File : BG051687.D
Acq On : 20 Dec 2021 18:48
Operator : CG/JU
Sample : PB141447BL
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK447

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/21/2021
Supervised By :mohammad ahmed 12/28/2021

Quant Time: Dec 20 23:11:47 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 14 14:19:57 2021
Response via : Initial Calibration



TIC: BG051687.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.596min (-0.009) 5.09 ng/uL****response 4690**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	71.34
34.00	0.00	0.00
0.00	0.00	0.00

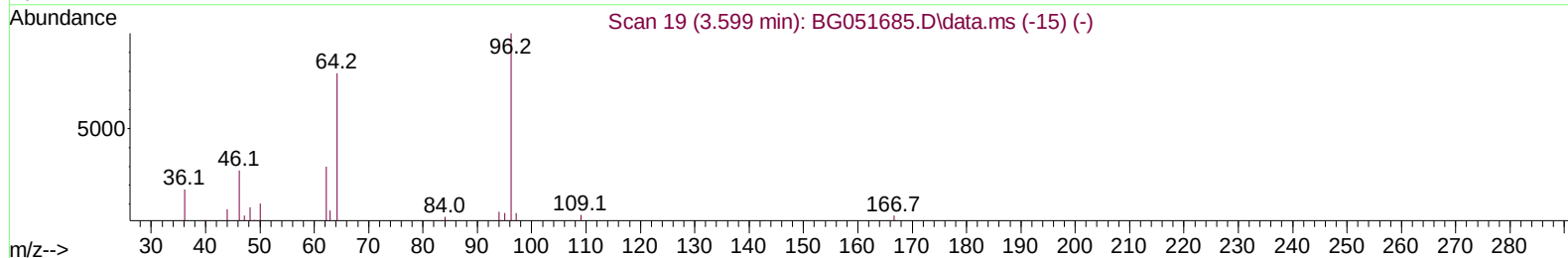
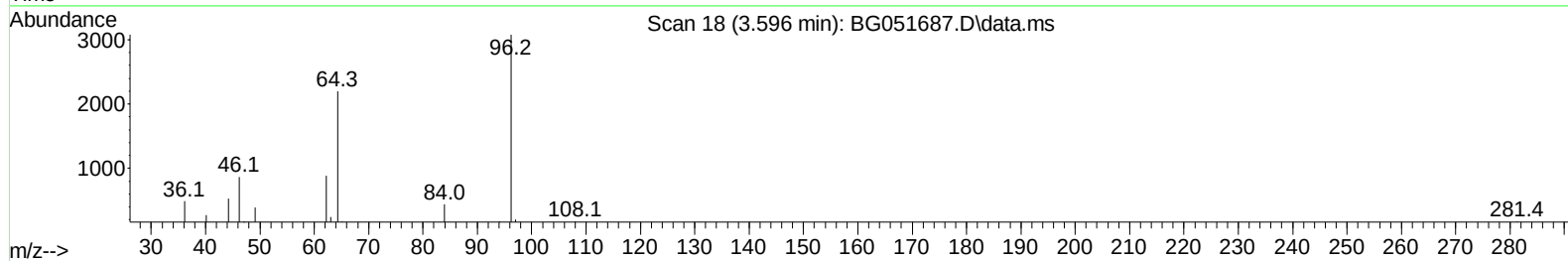
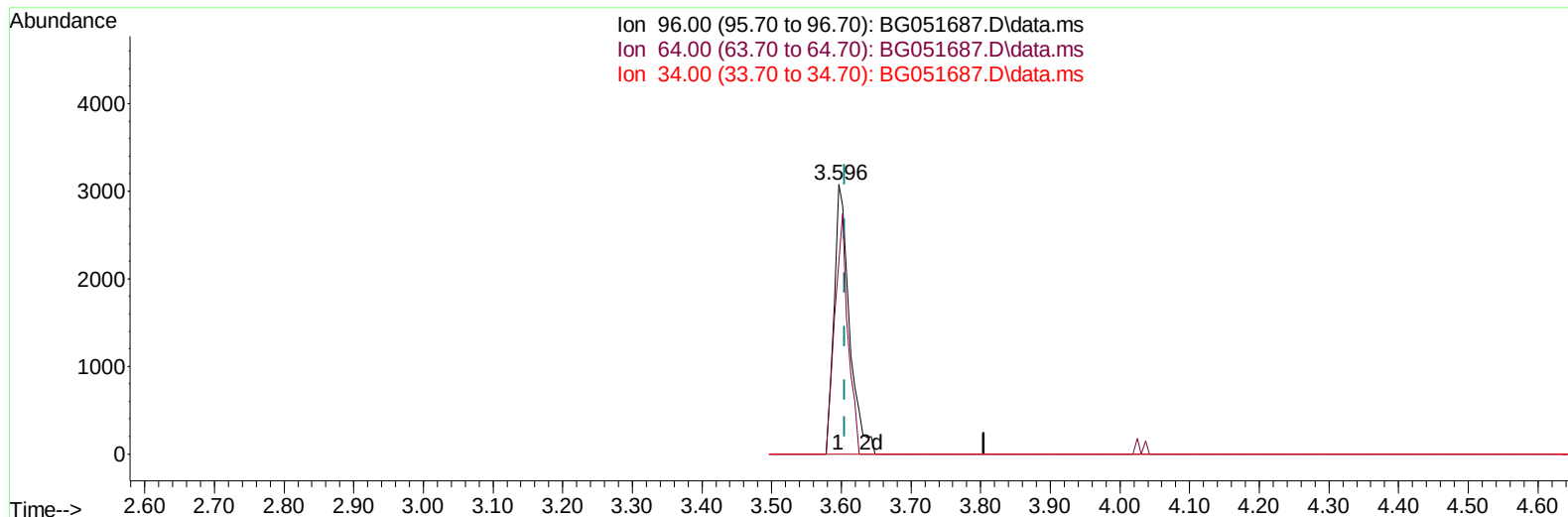
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(3) 1,4-Dioxane-d8 (S)

3.596min (-0.009) 5.16 ng/uL m

response 4761

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	71.34
34.00	0.00	0.00
0.00	0.00	0.00

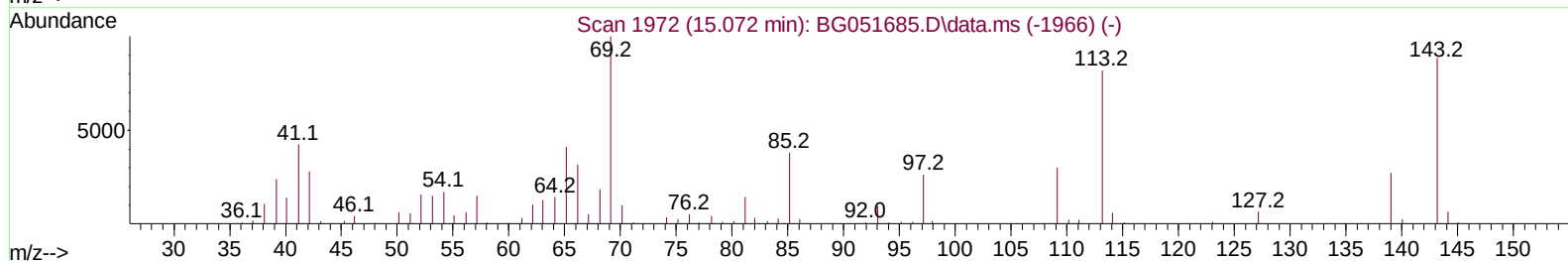
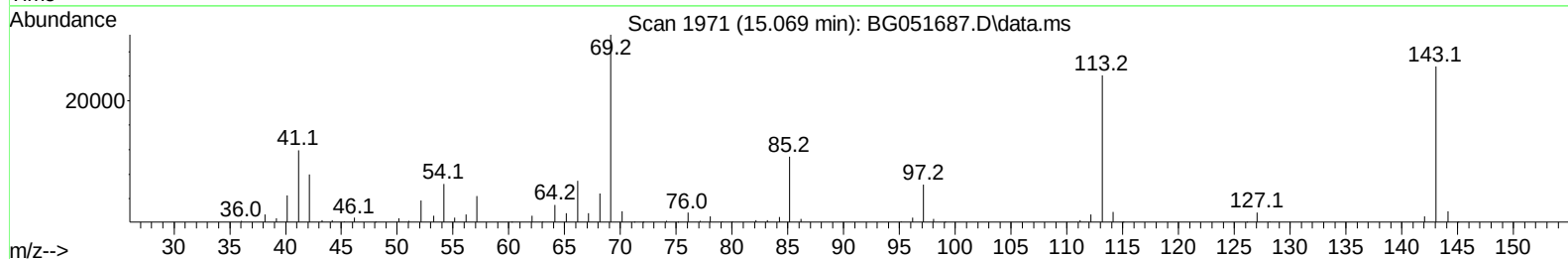
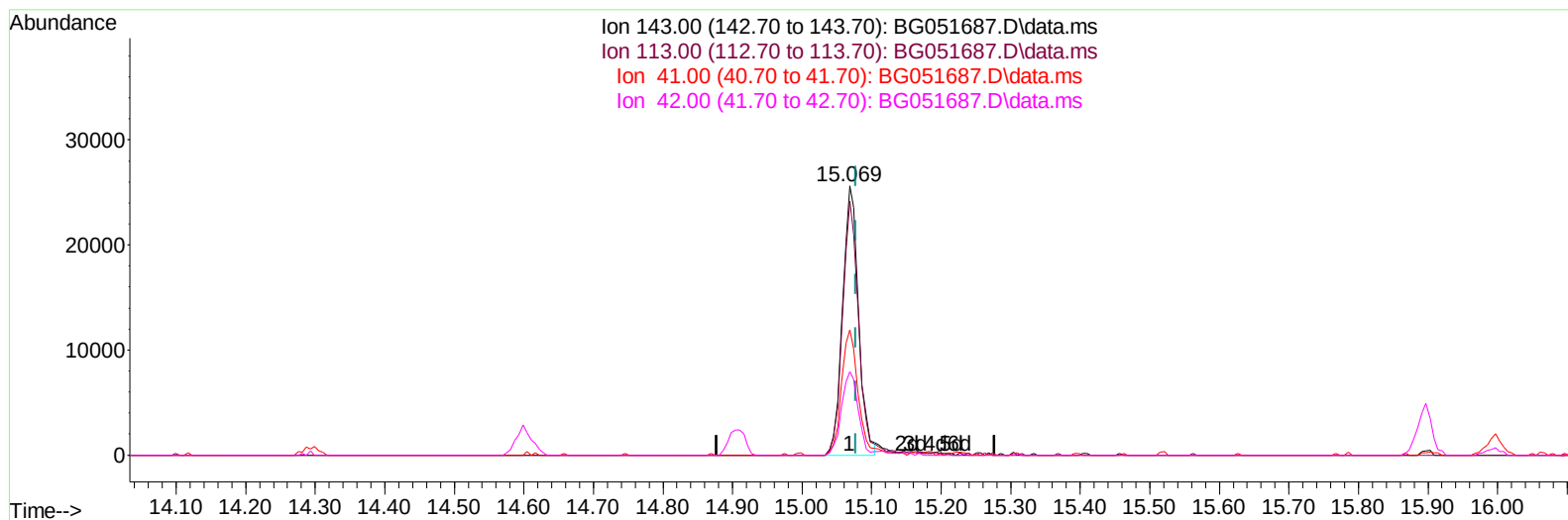
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TIC: BG051687.D\data.ms

(54) 4-Nitrophenol-d4 (S)**15.069min (-0.009) 27.25 ng/u1****response 41158**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.37
41.00	44.40	46.50
42.00	29.70	31.07

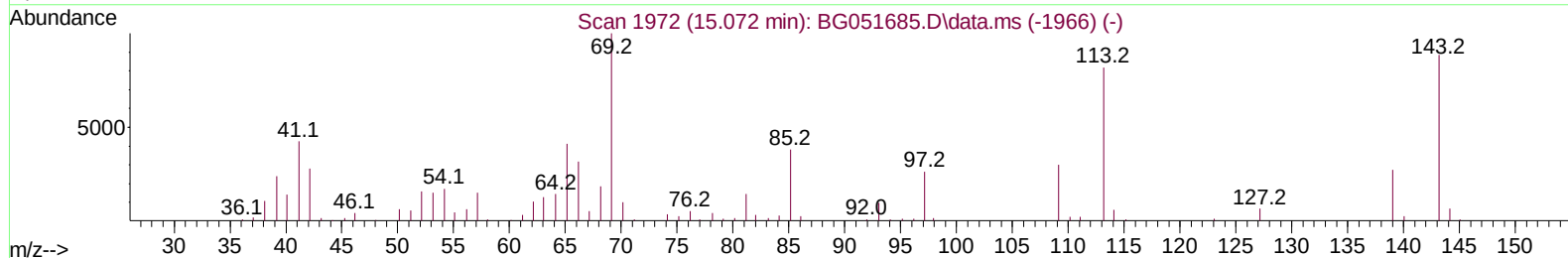
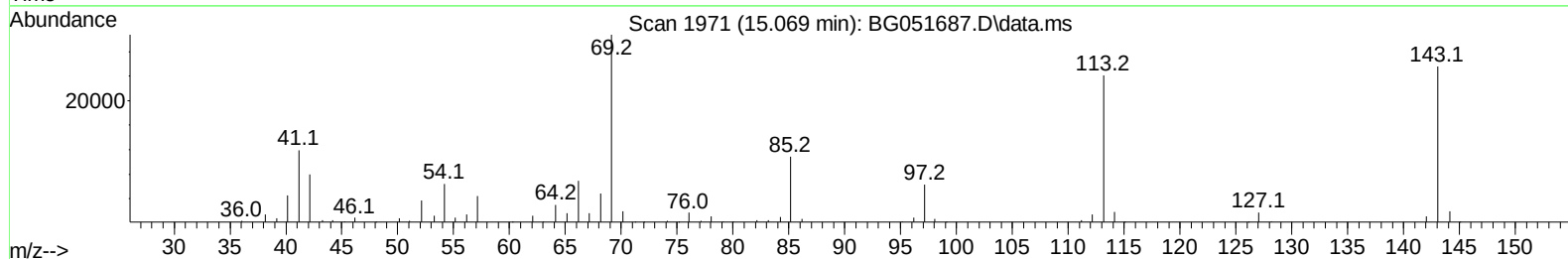
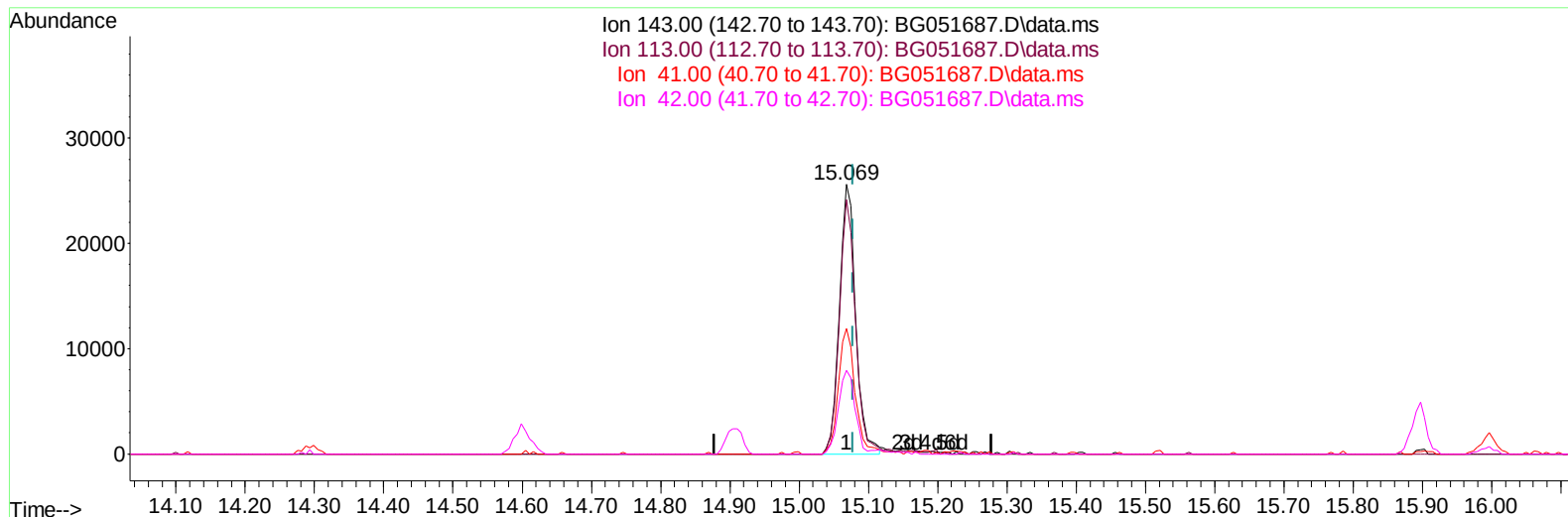
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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TIC: BG051687.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.069min (-0.009) 27.64 ng/ul m

response 41740

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	94.38
41.00	44.40	46.50
42.00	29.70	31.07

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.272	152	38290	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.109	136	164264	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.904	164	116037	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.653	188	269819	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.971	240	236538	20.000	ng/ul	#-0.01
88) Perylene-d12	25.466	264	234477	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	4761m	5.164	ng/uL	0.00
4) Pyridine-d5	4.025	84	70765	25.403	ng/ul	-0.02
7) Phenol-d5	7.403	99	107941	31.573	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.573	67	62142	30.576	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.796	132	74638	31.390	ng/ul	-0.01
15) 4-Methylphenol-d8	8.965	113	84646	32.031	ng/ul	-0.01
21) Nitrobenzene-d5	9.441	128	39759	33.014	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.169	143	44995	34.177	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.716	165	86800	30.403	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.227	131	97137	25.842	ng/ul	-0.01
46) Dimethylphthalate-d6	14.299	166	284736	30.658	ng/ul	-0.01
49) Acenaphthylene-d8	14.605	160	356470	30.904	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	41740m	27.636	ng/ul	0.00
60) Fluorene-d10	15.897	176	251876	30.270	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.997	200	30481	21.683	ng/ul	-0.01
73) Anthracene-d10	17.753	188	405292	31.488	ng/ul	0.00
81) Pyrene-d10	20.027	212	454930	31.956	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.231	264	399033	32.344	ng/ul	-0.02

Target Compounds Qvalue

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

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