

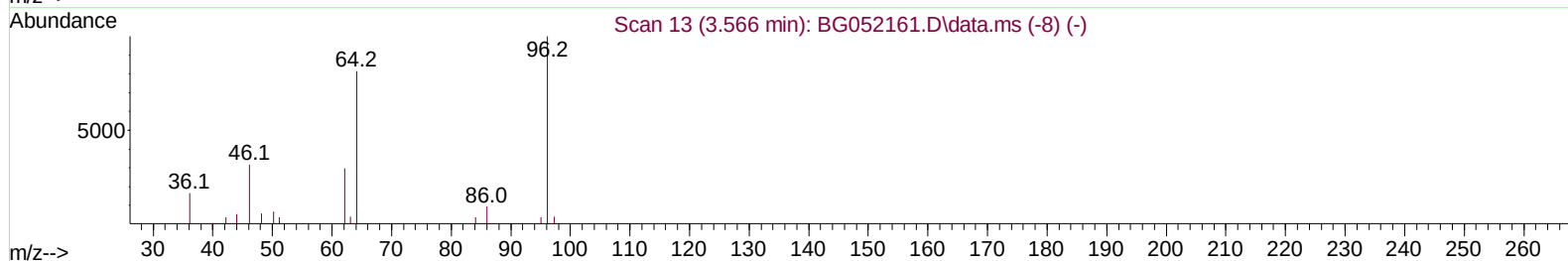
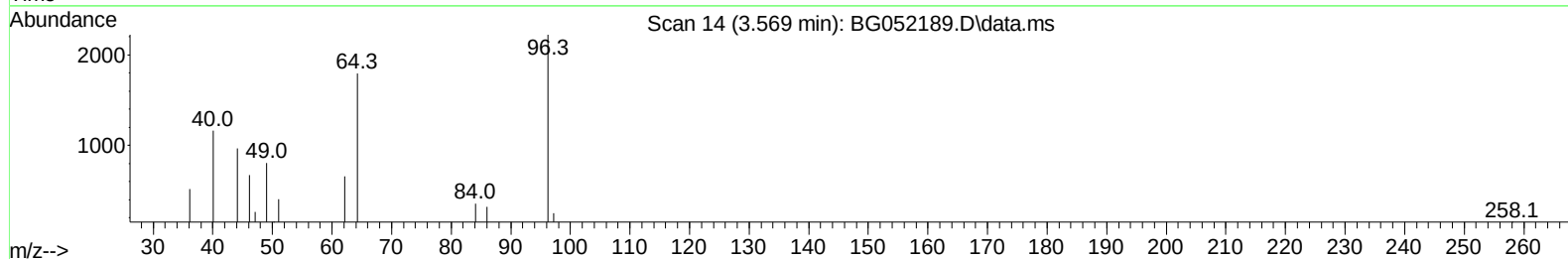
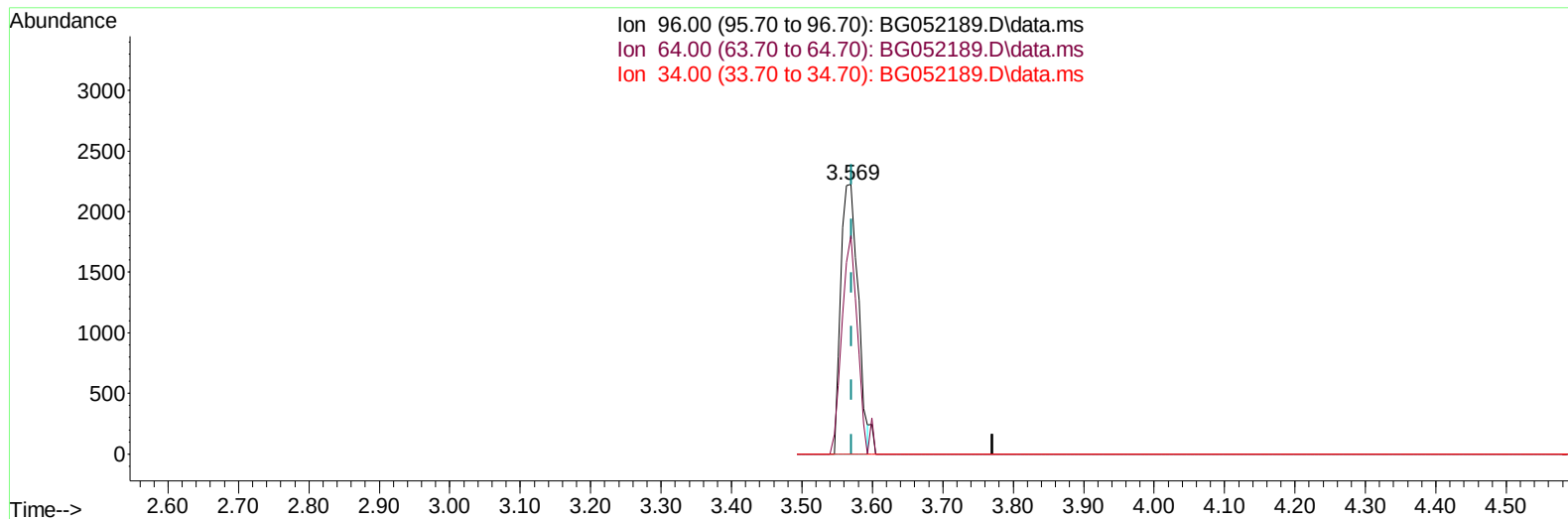
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052189.D
Acq On : 27 Jan 2022 17:20
Operator : CG/JU
Sample : N1273-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0HK0

Manual IntegrationsAPPROVED

Quant Time: Jan 28 00:00:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 26 23:58:52 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2022
Supervised By :Yogesh Patel 01/28/2022



TIC: BG052189.D\data.ms

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

3.569min (-0.001) 4.96 ng/uL

response 3744

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	80.89
34.00	0.00	0.00
0.00	0.00	0.00

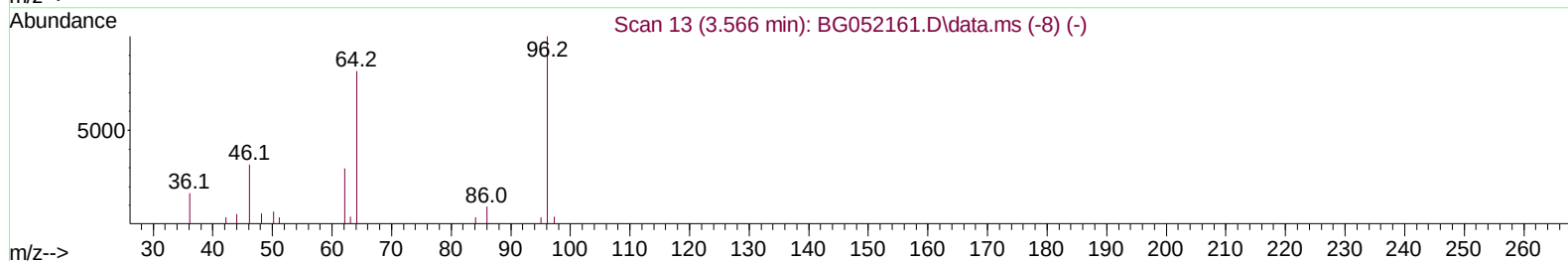
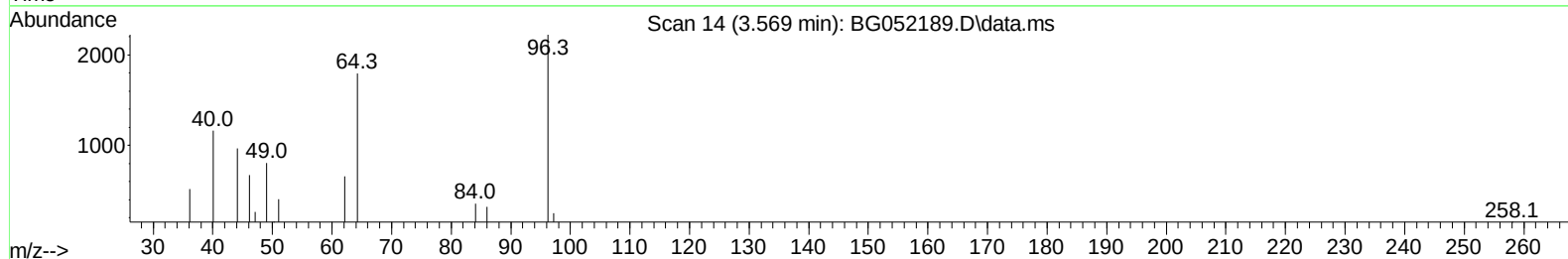
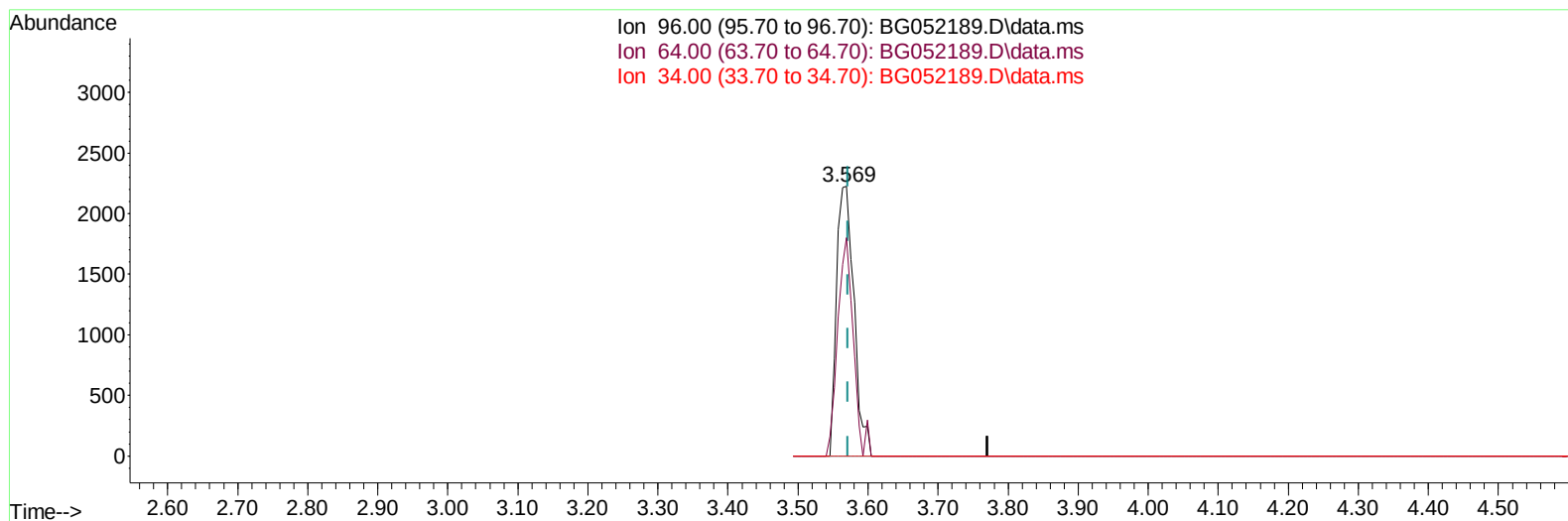
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TIC: BG052189.D\data.ms

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

3.569min (-0.001) 5.08 ng/uL m

response 3831

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	80.89
34.00	0.00	0.00
0.00	0.00	0.00

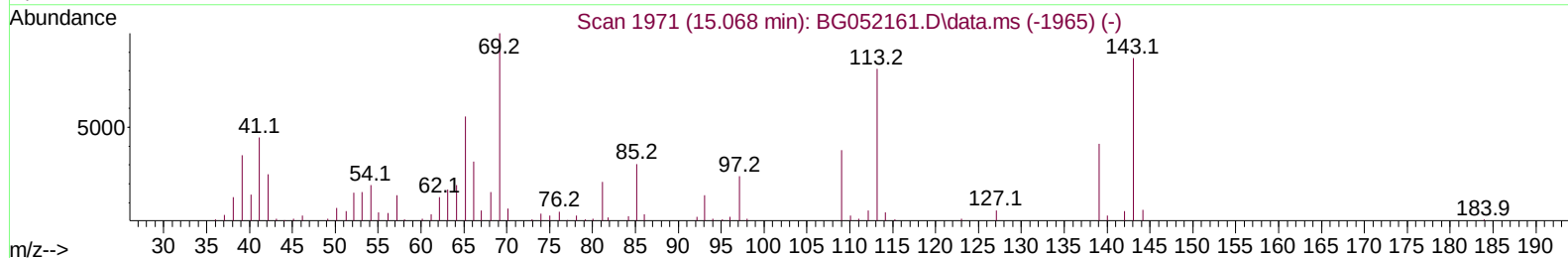
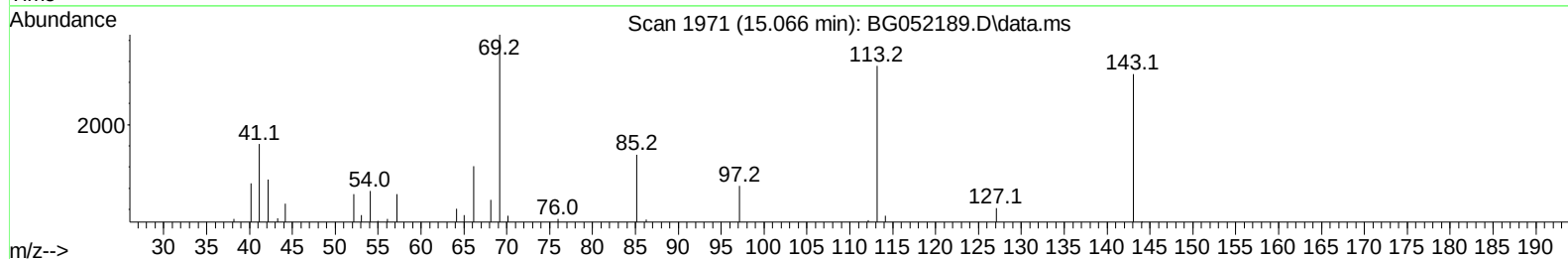
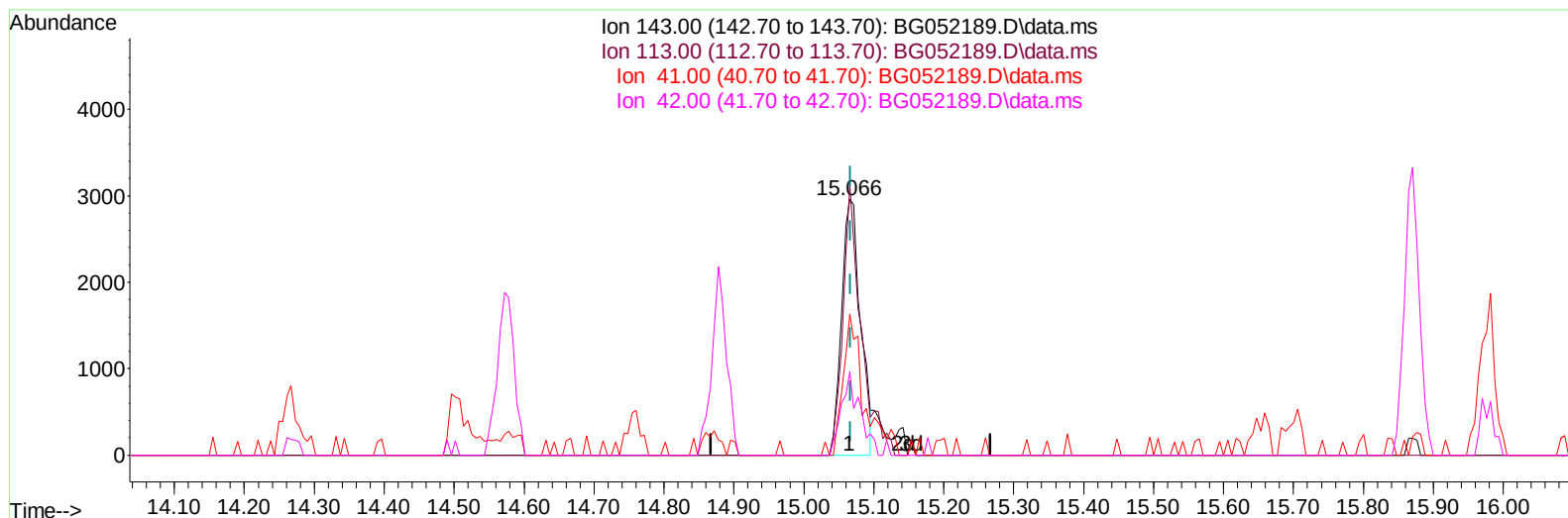
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Manual IntegrationsAPPROVED

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TIC: BG052189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (-0.001) 4.71 ng/u1

response 5625

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	105.17#
41.00	44.40	55.09#
42.00	29.70	32.51

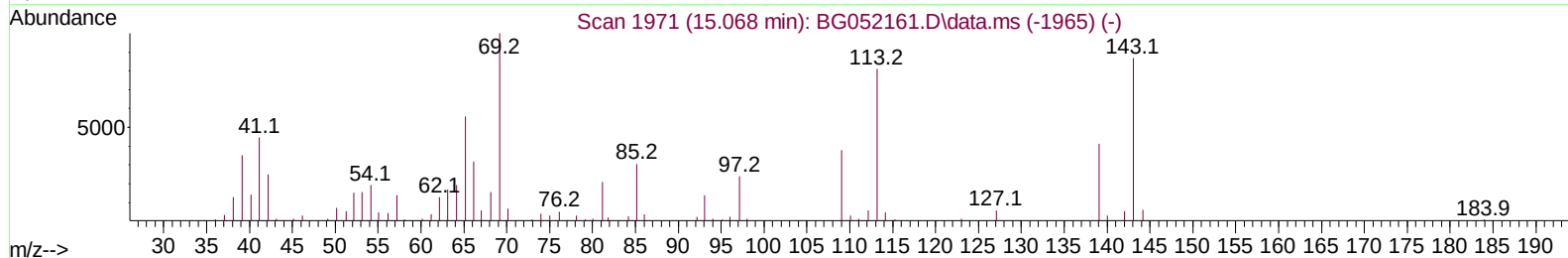
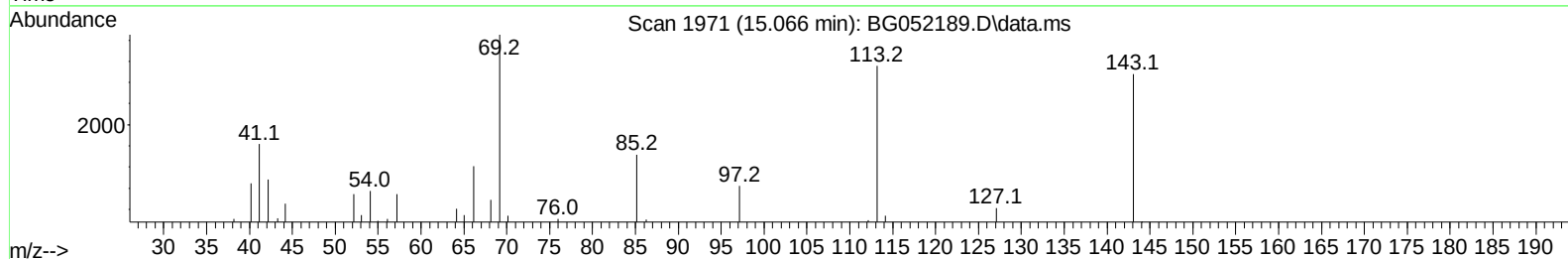
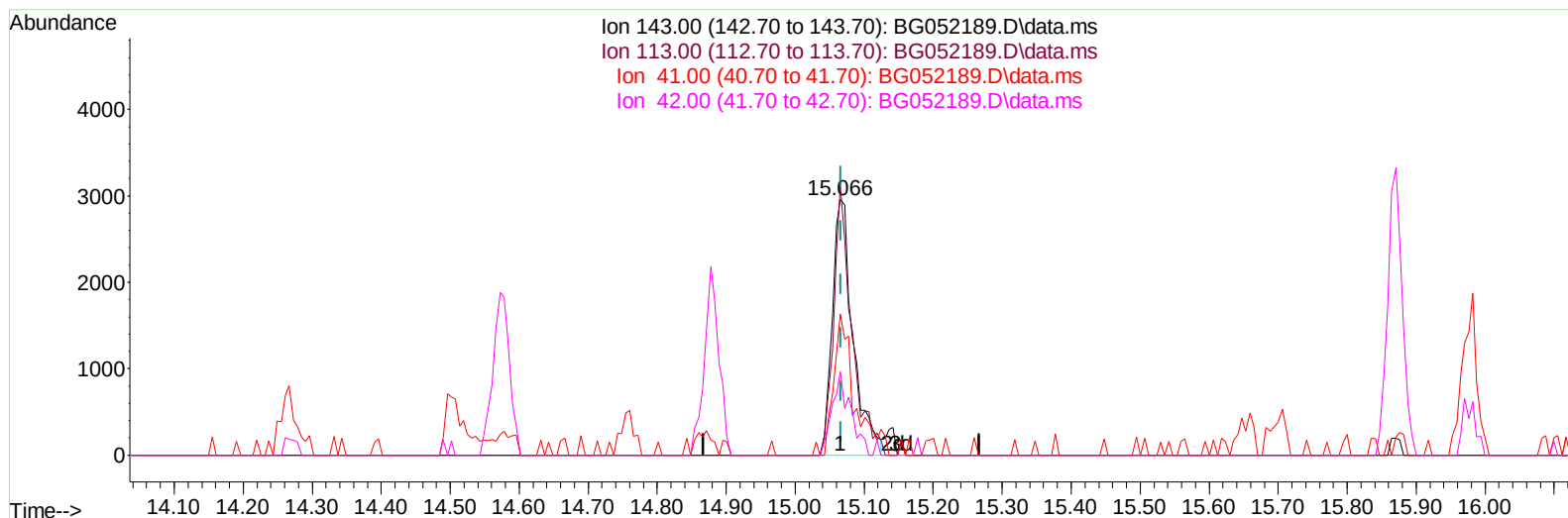
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Sample : N1273-01
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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TIC: BG052189.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.066min (-0.001) 5.20 ng/ul m

response 6202

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	105.17#
41.00	44.40	55.09#
42.00	29.70	32.51

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.240	152	26949	20.000	ng/ul	0.00
20) Naphthalene-d8	11.077	136	113032	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.884	164	86365	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	203738	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.945	240	199880	20.000	ng/ul	# 0.00
88) Perylene-d12	25.422	264	205278	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.569	96	3831m	5.080	ng/uL	0.00
4) Pyridine-d5	3.998	84	19791	8.442	ng/ul	0.00
7) Phenol-d5	7.382	99	16308	6.183	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.546	67	46493	27.768	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	40844	23.389	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	29650	14.461	ng/ul	0.00
21) Nitrobenzene-d5	9.409	128	28028	31.381	ng/ul	0.00
24) 2-Nitrophenol-d4	10.143	143	30446	30.822	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.689	165	56314	29.166	ng/ul	0.00
31) 4-Chloroaniline-d4	11.200	131	63378	23.570	ng/ul	0.00
46) Dimethylphthalate-d6	14.273	166	210900	29.543	ng/ul	0.00
49) Acenaphthylene-d8	14.578	160	262652	30.459	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143	6202m	5.196	ng/ul	0.00
60) Fluorene-d10	15.870	176	192451	30.920	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	28602	22.254	ng/ul	0.00
73) Anthracene-d10	17.727	188	328303	33.969	ng/ul	0.00
81) Pyrene-d10	20.006	212	393184	34.180	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.181	264	388627	35.962	ng/ul	0.00

Target Compounds Qvalue

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

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