

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
 Data File : BG043443.D
 Acq On : 31 Oct 2019 20:59
 Operator : JU
 Sample : K5615-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RBR-29

Quant Time: Nov 01 05:21:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 16:37:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	46902	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	164667	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	118109	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	265812	20.00	ng	0.00
76) Chrysene-d12	21.66	240	275729	20.00	ng	0.00
87) Perylene-d12	24.86	264	333022	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	320583	125.25	ng	0.00
7) Phenol-d6	7.21	99	441540	119.90	ng	0.00
23) Nitrobenzene-d5	9.18	82	265401	83.09	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	266450	118.55	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	717522	83.95	ng	0.00
79) Terphenyl-d14	20.01	244	1337489	91.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.10	163	43153	4.717	ng	98
71) Phenanthrene	17.44	178	39121	2.874	ng	98
75) Fluoranthene	19.45	202	40684	2.293	ng	95

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

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