

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
 Data File : BG046283.D
 Acq On : 14 Aug 2020 20:01
 Operator : CG/JU
 Sample : L3662-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RBR-59

Quant Time: Aug 17 01:31:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 18:43:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	90605	20.00	ng	0.00
21) Naphthalene-d8	10.74	136	382664	20.00	ng	0.00
39) Acenaphthene-d10	14.58	164	250213	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	526260	20.00	ng	0.00
76) Chrysene-d12	21.56	240	575842	20.00	ng	-0.01
86) Perylene-d12	24.67	264	661139	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	428806	82.27	ng	0.00
7) Phenol-d6	7.11	99	566925	79.81	ng	0.00
23) Nitrobenzene-d5	9.10	82	380161	53.79	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	295318	76.84	ng	0.00
45) 2-Fluorobiphenyl	13.20	172	916724	53.22	ng	0.00
79) Terphenyl-d14	19.93	244	1245130	44.04	ng	0.00
Target Compounds						
10) Phenol	7.14	94	16374	2.294	ng	83
20) 3+4-Methylphenols	8.71	107	110240	16.716	ng	90
50) Dimethylphthalate	14.02	163	117575	6.122	ng	# 95

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

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