

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
 Data File : BG041866.D
 Acq On : 1 Aug 2019 16:52
 Operator : HP/JU
 Sample : K4048-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWL-C6-GW

Quant Time: Aug 02 02:27:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 08:57:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	83623	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	300915	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	199191	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	473930	20.00	ng	0.00
76) Chrysene-d12	21.73	240	627817	20.00	ng	0.00
87) Perylene-d12	25.01	264	734517	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	254981	59.33	ng	0.00
7) Phenol-d6	7.19	99	215579	37.20	ng	0.00
23) Nitrobenzene-d5	9.21	82	455576	104.19	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	341466	150.92	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	1100102	97.41	ng	0.00
79) Terphenyl-d14	20.06	244	2114900	77.08	ng	0.00
Target Compounds						
10) Phenol	7.21	94	32242	5.349	ng	96
26) 2-Nitrophenol	9.96	139	4947	2.314	ng	# 89
32) Benzoic acid	10.12	122	43104	21.811	ng	93
50) Dimethylphthalate	14.14	163	125319	9.212	ng	99

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

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