

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
 Data File : BG041870.D
 Acq On : 1 Aug 2019 19:25
 Operator : HP/JU
 Sample : K4048-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWL-C1-GW

Quant Time: Aug 02 03:35:21 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	78114	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	289711	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	203786	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	484158	20.00	ng	0.00
76) Chrysene-d12	21.73	240	619990	20.00	ng	0.00
87) Perylene-d12	25.00	264	726121	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	294761	73.42	ng	0.00
7) Phenol-d6	7.19	99	271735	50.20	ng	0.00
23) Nitrobenzene-d5	9.20	82	447867	106.38	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	396255	171.19	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	1163727	100.72	ng	0.00
79) Terphenyl-d14	20.07	244	2237018	82.56	ng	0.00
Target Compounds						
10) Phenol	7.22	94	20534	3.647	ng	95
50) Dimethylphthalate	14.15	163	106917	7.682	ng	99
52) Acenaphthene	14.76	154	39185	3.611	ng	96
58) Fluorene	15.74	166	60507	4.543	ng	99

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

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