

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111419\
 Data File : BG043646.D
 Acq On : 14 Nov 2019 23:02
 Operator : JU
 Sample : K5670-05 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-03-111219

Quant Time: Nov 15 03:17:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	40102	20.00	ng	-0.01
21) Naphthalene-d8	10.80	136	156094	20.00	ng	-0.01
39) Acenaphthene-d10	14.63	164	105632	20.00	ng	-0.01
64) Phenanthrene-d10	17.37	188	234779	20.00	ng	-0.02
76) Chrysene-d12	21.64	240	256828	20.00	ng	-0.02
87) Perylene-d12	24.82	264	303573	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	74593	34.09	ng	0.00
7) Phenol-d6	7.20	99	101907	32.37	ng	0.00
23) Nitrobenzene-d5	9.17	82	60750	20.06	ng	0.00
42) 2,4,6-Tribromophenol	16.12	330	52282	26.01	ng	-0.01
45) 2-Fluorobiphenyl	13.25	172	167283	21.88	ng	-0.02
79) Terphenyl-d14	19.98	244	307073	22.43	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.08	163	18967	2.318	ng	Qvalue # 95
71) Phenanthrene	17.42	178	47947	3.988	ng	96
75) Fluoranthene	19.43	202	77229	4.927	ng	98
78) Pyrene	19.79	202	66490	4.182	ng	97
81) Benzo(a)anthracene	21.62	228	41395	2.573	ng	98
83) Chrysene	21.68	228	33556	2.099	ng	96
88) Benzo(b)fluoranthene	23.82	252	42881	2.494	ng	# 89
90) Benzo(a)pyrene	24.68	252	34448	2.098	ng	# 88

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

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