

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060319\
 Data File : BG041048.D
 Acq On : 3 Jun 2019 23:51
 Operator : HP/JU
 Sample : K3052-01 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FK-SF-03-007-A

Manual Integrations
 APPROVED

mohammad
 6/5/2019 8:35:33 AM

Quant Time: Jun 04 10:00:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	115732	20.00	ng	0.00
21) Naphthalene-d8	10.94	136	408029	20.00	ng	0.00
39) Acenaphthene-d10	14.75	164	259820	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	577396	20.00	ng	0.00
76) Chrysene-d12	21.78	240	596518	20.00	ng	0.00
87) Perylene-d12	25.13	264	560187	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	85553	13.33	ng	0.00
7) Phenol-d6	7.27	99	285583	28.81	ng	0.00
23) Nitrobenzene-d5	9.30	82	218044	23.72	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	13607	3.53	ng	-0.01
45) 2-Fluorobiphenyl	13.38	172	426231	23.62	ng	0.00
79) Terphenyl-d14	20.09	244	605875	21.56	ng	0.00
Target Compounds						
71) Phenanthrene	17.54	178	239201	7.953	ng	99
75) Fluoranthene	19.54	202	301478	8.115	ng	100
78) Pyrene	19.90	202	287014	7.402	ng	98
81) Benzo(a)anthracene	21.76	228	135970	3.424	ng	# 93
83) Chrysene	21.82	228	155187	4.084	ng	# 88
88) Benzo(b)fluoranthene	24.05	252	139562	4.108	ng	# 46
90) Benzo(a)pyrene	24.97	252	99260	3.062	ng	# 71

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

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