

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060916\
 Data File : BG022261.D
 Acq On : 9 Jun 2016 9:55
 Operator : UM/SJ
 Sample : H3402-07 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS-42

Quant Time: Jun 09 17:17:37 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 06 16:12:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	41459	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	213272	20.00	ng	0.00
38) Acenaphthene-d10	14.74	164	126260	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	248558	20.00	ng	0.00
75) Chrysene-d12	21.76	240	212662	20.00	ng	0.00
86) Perylene-d12	25.06	264	205729	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	112379	52.41	ng	0.00
7) Phenol-d6	7.27	99	176470	52.34	ng	0.00
23) Nitrobenzene-d5	9.28	82	92530	30.25	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	56761	39.79	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	265959	32.10	ng	0.00
78) Terphenyl-d14	20.08	244	295319	32.97	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	17.53	178	208363	15.43	ng	97
71) Anthracene	17.62	178	53182	3.97	ng	97
74) Fluoranthene	19.54	202	311587	20.08	ng	98
77) Pyrene	19.89	202	280727	21.23	ng	99
80) Benzo(a)anthracene	21.75	228	145511	12.20	ng	98
82) Chrysene	21.81	228	123352	10.63	ng	97
83) Bis(2-ethylhexyl)phthalate	21.62	149	21817	2.42	ng	94
85) Indeno(1,2,3-cd)pyrene	28.81	276	78080m	6.00	ng	
87) Benzo(b)fluoranthene	24.01	252	159431	13.29	ng	# 95
88) Benzo(k)fluoranthene	24.06	252	60467m	5.12	ng	
89) Benzo(a)pyrene	24.91	252	118416	10.32	ng	98
91) Benzo(a,h,i)perylene	30.01	276	69916m	6.22	ng	

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

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