

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
 Data File : BG017542.D
 Acq On : 8 Jun 2015 14:34
 Operator : TP/IZ
 Sample : G2464-18 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-16DS

Manual Integrations
 APPROVED

apatel
 6/9/2015 3:56:03 PM

Quant Time: Jun 09 02:24:54 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 09 02:10:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	16651	20.00	ng	0.00
21) Naphthalene-d8	10.41	136	69261	20.00	ng	0.00
38) Acenaphthene-d10	14.28	164	44179	20.00	ng	0.00
63) Phenanthrene-d10	17.03	188	96421	20.00	ng	0.00
75) Chrysene-d12	21.23	240	106878	20.00	ng	0.00
86) Perylene-d12	23.47	264	112688	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	16910	17.89	ng	0.00
7) Phenol-d6	6.82	99	24265	18.85	ng	0.00
23) Nitrobenzene-d5	8.78	82	18188	13.11	ng	0.00
41) 2,4,6-Tribromophenol	15.79	330	8102	13.20	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	40843	13.95	ng	0.00
78) Terphenyl-d14	19.67	244	64985	12.57	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	17.07	178	49226	9.16	ng	97
71) Anthracene	17.17	178	17126m	3.19	ng	
74) Fluoranthene	19.10	202	85148	12.11	ng	98
77) Pyrene	19.47	202	87131	13.23	ng	97
80) Benzo(a)anthracene	21.22	228	39857	6.06	ng	99
82) Chrysene	21.26	228	33097m	5.56	ng	
85) Indeno(1,2,3-cd)pyrene	25.73	276	20650	2.76	ng	# 93
87) Benzo(b)fluoranthene	22.79	252	51931m	7.27	ng	
88) Benzo(k)fluoranthene	22.83	252	14150m	2.07	ng	
89) Benzo(a)pyrene	23.36	252	37587	5.74	ng	# 97
91) Benzo(g,h,i)perylene	26.44	276	22021	3.40	ng	97

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

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