

Data Path : S:\HPCHEM1\BNA_M\DATA\BM041015\
 Data File : BM000947.D
 Acq On : 10 Apr 2015 19:49
 Operator : TP/IZ
 Sample : G1757-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LS-SW05

Manual Integrations
 APPROVED

apatel
 4/13/2015 4:23:57 PM

Quant Time: Apr 13 15:55:45 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM040915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 10 23:03:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	26607	5.00	ng	0.00
8) Naphthalene-d8	10.38	136	97418	5.00	ng	0.00
17) Acenaphthene-d10	14.25	164	42606	5.00	ng	-0.01
24) Phenanthrene-d10	16.99	188	99270	5.00	ng	-0.01
30) Chrysene-d12	21.20	240	87155	5.00	ng	0.00
38) Perylene-d12	23.39	264	80364	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	23718	4.21	ng	0.00
5) Phenol-d6	6.78	99	17915	2.66	ng	0.00
9) Nitrobenzene-d5	8.76	82	37400	7.95	ng	0.00
18) 2,4,6-Tribromophenol	15.75	330	23783	13.38	ng	-0.01
19) 2-Fluorobiphenyl	12.88	172	118991	9.97	ng	0.00
33) Terphenyl-d14	19.66	244	112547	10.61	ng	0.01
Target Compounds						Qvalue
13) Naphthalene	10.44	128	1077	0.05	ng	# 77
16) 1-Methylnaphthalene	12.29	142	315	0.02	ng	94
21) Acenaphthene	14.32	154	454m	0.04	ng	
27) Phenanthrene	17.04	178	1041	0.04	ng	# 76
29) Fluoranthene	19.07	202	882	0.03	ng	81
32) Pyrene	19.43	202	855	0.03	ng	# 83
34) Benzo(a)anthracene	21.18	228	933	0.04	ng	# 82
36) Chrysene	21.23	228	548m	0.02	ng	
37) Indeno(1,2,3-cd)pyrene	25.57	276	1309	0.05	ng	98
42) Dibenzo(a,h)anthracene	25.58	278	1389	0.07	ng	# 62
43) Benzo(g,h,i)perylene	26.23	276	849	0.04	ng	# 70

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

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