

Data Path : S:\HPCHEM1\BNA_M\DATA\BM041015\
 Data File : BM000942.D
 Acq On : 10 Apr 2015 16:50
 Operator : TP/IZ
 Sample : G1757-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LS-SW03

Manual Integrations
 APPROVED

apatel
 4/13/2015 4:23:37 PM

Quant Time: Apr 13 16:08:42 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 16:15:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	26199	5.00	ng	0.00
8) Naphthalene-d8	10.38	136	97307	5.00	ng	0.00
17) Acenaphthene-d10	14.25	164	42013	5.00	ng	0.00
24) Phenanthrene-d10	16.99	188	97960	5.00	ng	0.00
30) Chrysene-d12	21.20	240	84608	5.00	ng	0.00
38) Perylene-d12	23.39	264	76086	5.00	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	22753	4.10	ng	0.00
5) Phenol-d6	6.78	99	18533	2.79	ng	0.00
9) Nitrobenzene-d5	8.76	82	34367	7.31	ng	0.00
18) 2,4,6-Tribromophenol	15.75	330	23516	13.41	ng	0.00
19) 2-Fluorobiphenyl	12.88	172	119518	10.16	ng	0.00
33) Terphenyl-d14	19.66	244	114838	11.15	ng	0.02
Target Compounds						
27) Phenanthrene	17.04	178	567	0.02	ng	# 60
34) Benzo(a)anthracene	21.18	228	745	0.03	ng	# 80
37) Indeno(1,2,3-cd)pyrene	25.57	276	1647	0.06	ng	96
42) Dibenzo(a,h)anthracene	25.58	278	1856	0.10	ng	# 71
43) Benzo(g,h,i)perylene	26.23	276	1069	0.05	ng	# 73

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

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