

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040215\
 Data File : BM000808.D
 Acq On : 02 Apr 2015 13:13
 Operator : TP/IZ
 Sample : G1457-22DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RR-PAH-WPDL

Quant Time: Apr 03 02:11:06 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM033115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 04:17:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	27788	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	95026	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	46951	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	104693	5.00	ng	0.00
19) Chrysene-d12	21.21	240	87968	5.00	ng	0.00
25) Perylene-d12	23.40	264	80121	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	11104	2.05	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	38543	2.65	ng	0.00
21) Terphenyl-d14	19.66	244	29119	2.75	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	13761	0.62	ng	99
10) Acenaphthylene	13.98	152	14491	0.70	ng	99
11) Acenaphthene	14.33	154	50494	3.51	ng	99
12) Fluorene	15.32	166	31652	1.96	ng	100
16) Phenanthrene	17.05	178	21588	0.75	ng	99
17) Anthracene	17.15	178	10341m	0.43	ng	
18) Fluoranthene	19.08	202	26641	0.94	ng	100
20) Pyrene	19.44	202	18280	0.64	ng	99
22) Benzo(a)anthracene	21.20	228	6693	0.27	ng	99
23) Chrysene	21.25	228	12621	0.48	ng	# 90
24) Indeno(1,2,3-cd)pyrene	25.59	276	19330	0.71	ng	100
26) Benzo(b)fluoranthene	22.74	252	10431	0.41	ng	97
29) Dibenzo(a,h)anthracene	25.60	278	15372	0.72	ng	99
30) Benzo(g,h,i)perylene	26.26	276	19375	0.84	ng	99

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

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