

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040215\
 Data File : BM000807.D
 Acq On : 02 Apr 2015 10:58
 Operator : TP/IZ
 Sample : G1457-22
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RR-PAH-WP

Quant Time: Apr 03 02:09:21 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	32454	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	112535	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	56531	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	127473	5.00	ng	0.00
19) Chrysene-d12	21.21	240	102345	5.00	ng	0.00
25) Perylene-d12	23.40	264	93152	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	67631	9.75	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	199591	11.38	ng	0.00
21) Terphenyl-d14	19.66	244	157057	12.75	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	10.45	128	71079	2.72	ng	99
10) Acenaphthylene	13.98	152	82706	3.31	ng	100
11) Acenaphthene	14.33	154	265825	15.34	ng	98
12) Fluorene	15.32	166	171776	8.85	ng	99
16) Phenanthrene	17.05	178	116320	3.33	ng	99
17) Anthracene	17.13	178	54382m	1.86	ng	
18) Fluoranthene	19.08	202	147221	4.27	ng	100
20) Pyrene	19.44	202	100717	3.02	ng	99
22) Benzo(a)anthracene	21.19	228	36092	1.23	ng	# 91
23) Chrysene	21.24	228	64898	2.12	ng	96
24) Indeno(1,2,3-cd)pyrene	25.58	276	106295	3.35	ng	100
26) Benzo(b)fluoranthene	22.73	252	55372	1.89	ng	95
29) Dibenzo(a,h)anthracene	25.59	278	85240	3.43	ng	98
30) Benzo(g,h,i)perylene	26.25	276	103064	3.84	ng	99

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

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