

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020515\
 Data File : BE089514.D
 Acq On : 5 Feb 2015 22:09
 Operator : TP/IZ
 Sample : PB81591BSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81591BSD

Quant Time: Feb 06 04:12:42 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 06 03:21:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.49	152	27863	5.00	ng	0.00
2) Naphthalene-d8	12.40	136	115715	5.00	ng	0.00
6) Acenaphthene-d10	16.59	164	54569	5.00	ng	0.00
11) Phenanthrene-d10	19.90	188	104473	5.00	ng	0.00
16) Chrysene-d12	23.76	240	116856	5.00	ng	0.00
22) Perylene-d12	25.46	264	117520	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.76	82	61109	6.92	ng	0.00
7) 2-Fluorobiphenyl	15.03	172	112604	6.55	ng	0.00
18) Terphenyl-d14	22.40	244	148447	7.30	ng	0.00
Target Compounds						Ovalue
4) Naphthalene	12.44	128	148522	5.10	ng	99
5) 2-Methylnaphthalene	14.10	142	98721	5.17	ng	95
8) Acenaphthylene	16.24	152	643105	5.49	ng	100
9) Acenaphthene	16.65	154	94899	5.28	ng	100
10) Fluorene	17.91	166	115363	5.95	ng	98
12) Pentachlorophenol	19.58	266	15990	0.32	ng	94
13) Phenanthrene	19.94	178	180896	5.70	ng	99
14) Anthracene	20.04	178	181731	6.48	ng	100
15) Fluoranthene	21.83	202	187707	6.42	ng	99
17) Pyrene	22.15	202	201624	6.04	ng	99
19) Benzo(a)anthracene	23.75	228	774339	6.41	ng	98
20) Chrysene	23.79	228	771492	6.46	ng	99
21) Indeno(1,2,3-cd)pyrene	26.93	276	920201	6.96	ng	97
23) Benzo(b)fluoranthene	25.02	252	181077	6.11	ng	99
24) Benzo(k)fluoranthene	25.05	252	224184	5.98	ng	96
25) Benzo(a)pyrene	25.39	252	194686	6.64	ng	98
26) Dibenzo(a,h)anthracene	26.95	278	204525	6.76	ng	97
27) Benzo(a,h,i)perylene	27.36	276	784821	6.82	ng	95

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

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