

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020615\
 Data File : BE089543.D
 Acq On : 7 Feb 2015 2:54
 Operator : TP/IZ
 Sample : PB81635BSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81635BSD

Quant Time: Feb 07 06:02:29 2015
 Quant Method : Z:\HPCHEM1\BNA_E\methods\SIMPAH-BE020615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 07 04:07:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.42	152	29766	5.00	ng	0.00
2) Naphthalene-d8	12.33	136	105979	5.00	ng	0.00
6) Acenaphthene-d10	16.51	164	39382	5.00	ng	0.00
11) Phenanthrene-d10	19.84	188	67869	5.00	ng	0.00
16) Chrysene-d12	23.72	240	90075	5.00	ng	0.00
22) Perylene-d12	25.41	264	102324	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.69	82	61358	8.72	ng	0.00
7) 2-Fluorobiphenyl	14.96	172	86995	8.28	ng	0.00
18) Terphenyl-d14	22.36	244	89815	7.56	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	12.38	128	163045	6.83	ng	99
5) 2-Methylnaphthalene	14.03	142	91859	6.77	ng	98
8) Acenaphthylene	16.16	152	486985	7.35	ng	99
9) Acenaphthene	16.59	154	70949	6.74	ng	98
10) Fluorene	17.84	166	78859	7.07	ng	98
12) Pentachlorophenol	19.52	266	11891	12.33	ng	# 86
13) Phenanthrene	19.89	178	116492	6.93	ng	100
14) Anthracene	19.98	178	116071	8.17	ng	99
15) Fluoranthene	21.79	202	122526	7.82	ng	98
17) Pyrene	22.10	202	136430	7.11	ng	99
19) Benzo(a)anthracene	23.70	228	561650	7.01	ng	# 91
20) Chrysene	23.75	228	602199	7.40	ng	91
21) Indeno(1,2,3-cd)pyrene	26.85	276	863530	7.63	ng	99
23) Benzo(b)fluoranthene	24.98	252	141248	7.11	ng	98
24) Benzo(k)fluoranthene	25.01	252	194915	7.89	ng	98
25) Benzo(a)pyrene	25.35	252	155410	7.86	ng	98
26) Dibenzo(a,h)anthracene	26.89	278	191108	7.88	ng	# 96
27) Benzo(g,h,i)perylene	27.29	276	760970	7.69	ng	96

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

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