

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020615\
 Data File : BE089534.D
 Acq On : 6 Feb 2015 19:44
 Operator : TP/IZ
 Sample : PB81495BSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81495BSD

Quant Time: Feb 07 04:37:13 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	9.43	152	34826	5.00	ng	0.00
2) Naphthalene-d8	12.33	136	132735	5.00	ng	0.00
6) Acenaphthene-d10	16.51	164	55042	5.00	ng	0.00
11) Phenanthrene-d10	19.85	188	101331	5.00	ng	0.00
16) Chrysene-d12	23.72	240	120220	5.00	ng	0.00
22) Perylene-d12	25.41	264	126153	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.69	82	72941	8.28	ng	0.00
7) 2-Fluorobiphenyl	14.97	172	124362	8.47	ng	0.00
18) Terphenyl-d14	22.36	244	151722	9.57	ng	0.00

Target Compounds				Qvalue		
4) Naphthalene	12.38	128	184744	6.18	ng	99
5) 2-Methylnaphthalene	14.04	142	114776	6.76	ng	99
8) Acenaphthylene	16.16	152	704184	7.61	ng	99
9) Acenaphthene	16.59	154	103747	7.06	ng	98
10) Fluorene	17.84	166	122026	7.82	ng	99
12) Pentachlorophenol	19.52	266	24599	15.14	ng	88
13) Phenanthrene	19.89	178	189764	7.56	ng	99
14) Anthracene	19.99	178	189531	8.94	ng	100
15) Fluoranthene	21.79	202	197580	8.45	ng	99
17) Pyrene	22.10	202	214822	8.39	ng	99
19) Benzo(a)anthracene	23.71	228	843684	7.89	ng	97
20) Chrysene	23.75	228	869924	8.01	ng	97
21) Indeno(1,2,3-cd)pyrene	26.85	276	1114415	7.38	ng	99
23) Benzo(b)fluoranthene	24.98	252	214415	8.75	ng	98
24) Benzo(k)fluoranthene	25.01	252	251257	8.25	ng	99
25) Benzo(a)pyrene	25.35	252	226254	9.29	ng	98
26) Dibenzo(a,h)anthracene	26.89	278	248310	8.30	ng	97
27) Benzo(g,h,i)perylene	27.29	276	959168	7.86	ng	100

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

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