

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081815\
 Data File : BE090555.D
 Acq On : 18 Aug 2015 15:43
 Operator : UM/IZ
 Sample : PB85070BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB85070BS

Quant Time: Aug 19 02:31:36 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 19 02:30:21 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	23974	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	123302	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	67027	5.00	ng	0.00
18) Phenanthrene-d10	16.93	188	142875	5.00	ng	-0.02
25) Chrysene-d12	21.09	240	144000	5.00	ng	0.00
32) Perylene-d12	23.40	264	129604	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	41730	6.02	ng	0.00
5) Phenol-d6	6.77	99	64960	6.37	ng	0.00
7) Nitrobenzene-d5	8.72	82	29227	2.97	ng	0.00
13) 2,4,6-Tribromophenol	15.69	330	18284	7.09	ng	-0.02
14) 2-Fluorobiphenyl	12.81	172	58464	2.92	ng	0.00
27) Terphenyl-d14	19.55	244	85017	3.23	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.20	88	8165	2.98	ng	95
3) n-Nitrosodimethylamine	3.49	42	11123	2.69	ng	# 73
8) Nitrobenzene	8.76	77	29396	2.89	ng	93
9) Naphthalene	10.38	128	76154	2.72	ng	98
10) Hexachlorobutadiene	10.68	225	10875	2.64	ng	99
11) 2-Methylnaphthalene	12.00	142	51785	2.84	ng	97
15) Acenaphthylene	13.91	152	345193	2.82	ng	100
16) Acenaphthene	14.25	154	52182	2.83	ng	# 77
17) Fluorene	15.24	166	66971	2.97	ng	97
19) 4-Bromophenyl-phenylether	16.14	248	18078	3.10	ng	80
20) Hexachlorobenzene	16.25	284	79002	3.05	ng	96
21) Pentachlorophenol	16.60	266	18094	5.79	ng	91
22) Phenanthrene	16.98	178	111091	2.96	ng	100
23) Anthracene	17.07	178	112438	3.26	ng	100
24) Fluoranthene	18.98	202	129008	3.23	ng	99
26) Pyrene	19.34	202	136422	3.31	ng	99
28) Benzo(a)anthracene	21.08	228	120314	3.17	ng	99
29) Chrysene	21.13	228	122947	3.18	ng	97
30) Bis(2-ethylhexyl)phthalate	21.03	149	75643	3.39	ng	98
31) Indeno(1,2,3-cd)pyrene	25.77	276	117680	3.23	ng	# 92
33) Benzo(b)fluoranthene	22.69	252	107719	3.51	ng	98
34) Benzo(k)fluoranthene	22.74	252	114712	3.30	ng	98
35) Benzo(a)pyrene	23.29	252	101706	3.54	ng	98
36) Dibenzo(a,h)anthracene	25.79	278	93483	3.37	ng	97
37) Benzo(g,h,i)perylene	26.51	276	103365	3.41	ng	95

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

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