

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
 Data File : BM001529.D
 Acq On : 03 Jun 2015 00:49
 Operator : TP/IZ
 Sample : PB83718BSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB83718BSD

Quant Time: Jun 03 03:19:25 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 02:19:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	15960	5.00	ng	0.00
6) Naphthalene-d8	10.49	136	55085	5.00	ng	0.00
12) Acenaphthene-d10	14.36	164	25885	5.00	ng	-0.01
18) Phenanthrene-d10	17.10	188	67872	5.00	ng	0.00
25) Chrysene-d12	21.30	240	42652	5.00	ng	0.00
32) Perylene-d12	23.54	264	38326	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.28	112	48615	13.93	ng	-0.02
5) Phenol-d6	6.86	99	62274	14.44	ng	0.00
7) Nitrobenzene-d5	8.87	82	35998	10.50	ng	0.00
13) 2,4,6-Tribromophenol	15.84	330	15794	32.72	ng	0.00
14) 2-Fluorobiphenyl	12.98	172	76667	9.86	ng	-0.01
27) Terphenyl-d14	19.74	244	51756	9.84	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.23	88	13460	9.75	ng	98
3) n-Nitrosodimethylamine	3.52	42	22754	10.04	ng	93
8) Nitrobenzene	8.91	77	40710	11.18	ng	98
9) Naphthalene	10.54	128	108684	9.35	ng	97
10) Hexachlorobutadiene	10.83	225	19487	9.61	ng	100
11) 2-Methylnaphthalene	12.16	142	72210	9.90	ng	97
15) Acenaphthylene	14.08	152	114681	10.37	ng	100
16) Acenaphthene	14.42	154	64977	9.40	ng	99
17) Fluorene	15.40	166	61641	5.35	ng	93
19) 4-Bromophenyl-phenylether	16.28	248	20266	21.44	ng	# 58
20) Hexachlorobenzene	16.42	284	13087	5.15	ng	# 37
21) Pentachlorophenol	16.76	266	122563	81.84	ng	# 89
22) Phenanthrene	17.13	178	175883	14.82	ng	95
23) Anthracene	17.23	178	79199	5.52	ng	94
24) Fluoranthene	19.16	202	129450	10.87	ng	98
26) Pyrene	19.52	202	138375	10.75	ng	99
28) Benzo(a)anthracene	21.27	228	118423	10.31	ng	97
29) Chrysene	21.27	228	118423	10.89	ng	93
30) Bis(2-ethylhexyl)phthalate	21.20	149	97633	12.04	ng	98
31) Indeno(1,2,3-cd)pyrene	25.80	276	123058	10.18	ng	# 93
33) Benzo(b)fluoranthene	22.85	252	112341	10.27	ng	97
34) Benzo(k)fluoranthene	22.90	252	110488	11.55	ng	98
35) Benzo(a)pyrene	23.44	252	106058	11.37	ng	98
36) Dibenzo(a,h)anthracene	25.82	278	100156	10.79	ng	98
37) Benzo(g,h,i)perylene	26.50	276	103142	10.62	ng	96

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

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