

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
 Data File : BM001528.D
 Acq On : 03 Jun 2015 00:13
 Operator : TP/IZ
 Sample : PB83718BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB83718BS

Manual Integrations
 APPROVED

apatel
 6/3/2015 5:28:41 PM

Quant Time: Jun 03 03:18:39 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	14817	5.00	ng	0.00
6) Naphthalene-d8	10.49	136	50555	5.00	ng	0.00
12) Acenaphthene-d10	14.36	164	23648	5.00	ng	-0.01
18) Phenanthrene-d10	17.10	188	65703	5.00	ng	0.00
25) Chrysene-d12	21.30	240	41206	5.00	ng	0.00
32) Perylene-d12	23.53	264	36266	5.00	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.28	112	45465	14.04	ng	-0.02
5) Phenol-d6	6.86	99	57879	14.46	ng	0.00
7) Nitrobenzene-d5	8.87	82	33024	10.50	ng	0.00
13) 2,4,6-Tribromophenol	15.84	330	13846	31.40	ng	0.00
14) 2-Fluorobiphenyl	12.98	172	70528	9.93	ng	-0.01
27) Terphenyl-d14	19.74	244	49938	9.83	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	12723	9.93	ng	96
3) n-Nitrosodimethylamine	3.52	42	21471	10.20	ng	# 96
8) Nitrobenzene	8.91	77	36793	11.01	ng	98
9) Naphthalene	10.54	128	99281	9.30	ng	97
10) Hexachlorobutadiene	10.83	225	17777	9.55	ng	100
11) 2-Methylnaphthalene	12.16	142	64989	9.70	ng	98
15) Acenaphthylene	14.08	152	103807	10.28	ng	99
16) Acenaphthene	14.42	154	59005	9.34	ng	99
17) Fluorene	15.40	166	64539	6.13	ng	93
19) 4-Bromophenyl-phenylether	16.28	248	18332	20.03	ng	# 56
20) Hexachlorobenzene	16.42	284	12892	5.24	ng	# 37
21) Pentachlorophenol	16.76	266	117971	81.38	ng	# 89
22) Phenanthrene	17.13	178	163927	14.27	ng	96
23) Anthracene	17.23	178	80683	5.81	ng	95
24) Fluoranthene	19.16	202	122628	10.64	ng	98
26) Pyrene	19.52	202	132750	10.68	ng	99
28) Benzo(a)anthracene	21.26	228	113401	10.22	ng	96
29) Chrysene	21.33	228	107423m	10.22	ng	
30) Bis(2-ethylhexyl)phthalate	21.20	149	88576	11.31	ng	# 97
31) Indeno(1,2,3-cd)pyrene	25.80	276	113568	9.72	ng	# 94
33) Benzo(b)fluoranthene	22.85	252	105471	10.19	ng	97
34) Benzo(k)fluoranthene	22.90	252	104000	11.49	ng	99
35) Benzo(a)pyrene	23.44	252	99887	11.32	ng	98
36) Dibenzo(a,h)anthracene	25.82	278	92012	10.47	ng	97
37) Benzo(g,h,i)perylene	26.49	276	94640	10.30	ng	# 93

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

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