

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060815\
 Data File : BM001623.D
 Acq On : 09 Jun 2015 20:11
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Manual Integrations
 APPROVED

apatel
 6/10/2015 1:05:59 PM

Quant Time: Jun 10 04:04:22 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 10 03:26:36 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	20881	5.00	ng	0.00
6) Naphthalene-d8	10.47	136	75196	5.00	ng	0.00
12) Acenaphthene-d10	14.35	164	35917	5.00	ng	0.00
18) Phenanthrene-d10	17.09	188	44372	5.00	ng	0.00
25) Chrysene-d12	21.27	240	64096	5.00	ng	-0.02
32) Perylene-d12	23.51	264	58053	5.00	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.27	112	4758	1.00	ng	0.00
5) Phenol-d6	6.84	99	5909	1.01	ng	0.00
7) Nitrobenzene-d5	8.85	82	5257	0.99	ng	0.00
13) 2,4,6-Tribromophenol	15.83	330	1457	0.96	ng	0.00
14) 2-Fluorobiphenyl	12.96	172	11306	0.97	ng	0.00
27) Terphenyl-d14	19.72	244	8155	0.98	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.21	88	2147	0.93	ng	98
3) n-Nitrosodimethylamine	3.51	42	3219	0.98	ng	99
8) Nitrobenzene	8.89	77	5636	0.99	ng	100
9) Naphthalene	10.52	128	16792	1.00	ng	99
10) Hexachlorobutadiene	10.82	225	2807	0.97	ng	99
11) 2-Methylnaphthalene	12.15	142	10375	1.01	ng	96
15) Acenaphthylene	14.06	152	15834	0.98	ng	100
16) Acenaphthene	14.40	154	9886	0.97	ng	99
17) Fluorene	15.36	166	7739	1.13	ng	97
19) 4-Bromophenyl-phenylether	16.27	248	4441	1.06	ng	# 74
20) Hexachlorobenzene	16.38	284	2975	1.29	ng	# 80
21) Pentachlorophenol	16.72	266	954	1.24	ng	97
22) Phenanthrene	17.12	178	15048	0.95	ng	99
23) Anthracene	17.19	178	12793m	1.36	ng	
24) Fluoranthene	19.14	202	18766	1.09	ng	100
26) Pyrene	19.50	202	19855	0.96	ng	99
28) Benzo(a)anthracene	21.25	228	17971	0.96	ng	97
29) Chrysene	21.31	228	17260	0.98	ng	98
30) Bis(2-ethylhexyl)phthalate	21.19	149	9610	0.89	ng	99
31) Indeno(1,2,3-cd)pyrene	25.76	276	17751	0.99	ng	99
33) Benzo(b)fluoranthene	22.83	252	17452	1.00	ng	98
34) Benzo(k)fluoranthene	22.88	252	16155	0.96	ng	98
35) Benzo(a)pyrene	23.41	252	15109	0.98	ng	97
36) Dibenzo(a,h)anthracene	25.78	278	13891	0.99	ng	96
37) Benzo(g,h,i)perylene	26.45	276	14746	0.97	ng	98

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

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