

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
 Data File : BM001518.D
 Acq On : 02 Jun 2015 18:17
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Manual Integrations
 APPROVED

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 6/3/2015 5:28:10 PM

Quant Time: Jun 03 02:12:08 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 15:30:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	19804	5.00	ng	0.00
6) Naphthalene-d8	10.49	136	70118	5.00	ng	0.00
12) Acenaphthene-d10	14.36	164	33868	5.00	ng	0.00
18) Phenanthrene-d10	17.09	188	93452	5.00	ng	0.00
25) Chrysene-d12	21.30	240	54269	5.00	ng	0.00
32) Perylene-d12	23.54	264	48605	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.28	112	8909	2.04	ng	-0.02
5) Phenol-d6	6.86	99	10983	2.04	ng	0.00
7) Nitrobenzene-d5	8.87	82	8899	2.06	ng	0.00
13) 2,4,6-Tribromophenol	15.83	330	1202	1.65	ng	0.00
14) 2-Fluorobiphenyl	12.98	172	20512	2.04	ng	0.00
27) Terphenyl-d14	19.75	244	14094	2.08	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	3659	1.81	ng	98
3) n-Nitrosodimethylamine	3.52	42	5996	2.13	ng	98
8) Nitrobenzene	8.91	77	9572	2.08	ng	97
9) Naphthalene	10.54	128	29692	2.01	ng	98
10) Hexachlorobutadiene	10.83	225	5194	2.01	ng	100
11) 2-Methylnaphthalene	12.17	142	18909	2.01	ng	96
15) Acenaphthylene	14.07	152	29949	2.05	ng	99
16) Acenaphthene	14.42	154	18553	2.07	ng	95
17) Fluorene	15.39	166	30751	2.19	ng	98
19) 4-Bromophenyl-phenylether	16.31	248	2804	1.79	ng	# 43
20) Hexachlorobenzene	16.41	284	7378	2.26	ng	# 60
21) Pentachlorophenol	16.75	266	3434	2.37	ng	90
22) Phenanthrene	17.12	178	32521	1.82	ng	99
23) Anthracene	17.23	178	40216	2.16	ng	99
24) Fluoranthene	19.16	202	33174	1.95	ng	98
26) Pyrene	19.53	202	33982	2.05	ng	99
28) Benzo(a)anthracene	21.27	228	29181	1.97	ng	93
29) Chrysene	21.33	228	27775m	2.01	ng	
30) Bis(2-ethylhexyl)phthalate	21.22	149	20458	1.67	ng	# 99
31) Indeno(1,2,3-cd)pyrene	25.80	276	30654	2.04	ng	# 93
33) Benzo(b)fluoranthene	22.87	252	28616	2.03	ng	96
34) Benzo(k)fluoranthene	22.91	252	25019	1.99	ng	97
35) Benzo(a)pyrene	23.44	252	24488	2.02	ng	# 96
36) Dibenzo(a,h)anthracene	25.83	278	24380	2.10	ng	# 94
37) Benzo(g,h,i)perylene	26.50	276	25038	2.05	ng	# 94

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

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