Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060615\
Data File : BM001569.D
Acq On : 05 Jun 2015 17:14
Operator : TP/IZ
Sample : SSTDICC005
Misc :
ALS Vial : 1 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC005

Quant Time: Jun 06 01:19:08 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jun 06 01:16:45 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	33	5.00	ng		0.00
6) Naphthalene-d8	10.47	136	136	5.00	ng		0.00
12) Acenaphthene-d10	14.35	164	79	5.00	ng		0.00
18) Phenanthrene-d10	17.09	188	157	5.00	ng		0.00
25) Chrysene-d12	21.28	240	280	5.00	ng		0.00
32) Perylene-d12	23.52	264	341	5.00	ng		0.01
System Monitoring Compounds							
4) 2-Fluorophenol	5.28	112	82	11.03	ng		0.00
5) Phenol-d6	6.86	99	169	18.48	ng		0.00
7) Nitrobenzene-d5	8.86	82	107	11.17	ng		0.00
13) 2,4,6-Tribromophenol	15.83	330	37	11.28	ng		0.00
14) 2-Fluorobiphenyl	12.97	172	189	8.04	ng		0.00
27) Terphenyl-d14	19.72	244	157	4.53	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.21	88	108	31.27	ng	#	41
3) n-Nitrosodimethylamine	3.52	42	65	12.72	ng	#	90
8) Nitrobenzene	8.90	77	117	11.57	ng	#	91
9) Naphthalene	10.52	128	376	13.13	ng	#	59
10) Hexachlorobutadiene	10.82	225	60	12.15	ng		92
11) 2-Methylnaphthalene	12.15	142	253	14.08	ng	#	90
15) Acenaphthylene	14.06	152	372	10.87	ng	#	94
16) Acenaphthene	14.41	154	245	11.80	ng		99
17) Fluorene	15.39	166	179	10.98	ng	#	96
19) 4-Bromophenyl-phenylether	16.27	248	126	19.03	ng		96
20) Hexachlorobenzene	16.38	284	61	17.68	ng	#	69
21) Pentachlorophenol	16.75	266	42874	23537.51	ng		92
22) Phenanthrene	17.12	178	634	13.63	ng	#	79
23) Anthracene	17.12	178	862	40.99	ng	#	59
24) Fluoranthene	19.15	202	559	16.83	ng	#	90
26) Pyrene	19.52	202	660	7.64	ng		98
28) Benzo(a)anthracene	21.25	228	844	10.77	ng	#	81
29) Chrysene	21.31	228	797	11.35	ng	#	87
30) Bis(2-ethylhexyl)phthalate	21.19	149	860	15.50	ng	#	96
31) Indeno(1,2,3-cd)pyrene	25.77	276	2480	32.15	ng		98
33) Benzo(b)fluoranthene	22.84	252	1344	14.21	ng	#	51
34) Benzo(k)fluoranthene	22.89	252	1172	13.24	ng	#	69
35) Benzo(a)pyrene	23.42	252	1511	17.68	ng	#	54
36) Dibenzo(a,h)anthracene	25.79	278	2132	26.73	ng	#	83
37) Benzo(g,h,i)perylene	26.47	276	2002	23.87	ng	#	86

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

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