Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
Data File : BM001523.D
Acq On : 02 Jun 2015 21:15
Operator : TP/IZ
Sample : SSTDICC0.1
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.1

Quant Time: Jun 03 01:40:53 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	17154	5.00	ng		0.00
6) Naphthalene-d8	10.49	136	60663	5.00	ng		0.00
12) Acenaphthene-d10	14.36	164	29657	5.00	ng		0.00
18) Phenanthrene-d10	17.09	188	92700	5.00	ng		0.00
25) Chrysene-d12	21.30	240	54667	5.00	ng		0.00
32) Perylene-d12	23.54	264	50086	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.28	112	374	0.10	ng		-0.02
5) Phenol-d6	6.86	99	467	0.10	ng		0.00
7) Nitrobenzene-d5	8.87	82	367	0.10	ng		0.00
13) 2,4,6-Tribromophenol	15.83	330	55	0.09	ng		0.00
14) 2-Fluorobiphenyl	12.98	172	897	0.10	ng		0.00
27) Terphenyl-d14	19.75	244	669	0.10	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.23	88	276	0.16	ng	#	69
3) n-Nitrosodimethylamine	3.54	42	239	0.10	ng	#	89
8) Nitrobenzene	8.91	77	388	0.10	ng		97
9) Naphthalene	10.54	128	1326	0.10	ng	#	88
10) Hexachlorobutadiene	10.83	225	228	0.10	ng		98
11) 2-Methylnaphthalene	12.17	142	823	0.10	ng		95
15) Acenaphthylene	14.07	152	1210	0.09	ng		100
16) Acenaphthene	14.42	154	788	0.10	ng		98
17) Fluorene	15.39	166	1323	0.11	ng		99
19) 4-Bromophenyl-phenylether	16.31	248	118	0.08	ng	#	54
20) Hexachlorobenzene	16.41	284	342	0.11	ng	#	59
21) Pentachlorophenol	16.75	266	100	0.07	ng	#	85
22) Phenanthrene	17.12	178	1798	0.10	ng	#	95
23) Anthracene	17.23	178	1787	0.10	ng		100
24) Fluoranthene	19.16	202	1628	0.10	ng		98
26) Pyrene	19.53	202	1683	0.10	ng		100
28) Benzo(a)anthracene	21.27	228	1584	0.11	ng	#	91
29) Chrysene	21.33	228	1458	0.10	ng	#	87
30) Bis(2-ethylhexyl)phthalate	21.21	149	1185	0.10	ng	#	95
31) Indeno(1,2,3-cd)pyrene	25.81	276	1559	0.10	ng	#	93
33) Benzo(b)fluoranthene	22.85	252	1374	0.09	ng	#	63
34) Benzo(k)fluoranthene	22.91	252	1350	0.10	ng	#	59
35) Benzo(a)pyrene	23.44	252	1220	0.10	ng	#	53
36) Dibenzo(a,h)anthracene	25.83	278	1248	0.10	ng	#	70
37) Benzo(g,h,i)perylene	26.49	276	1321	0.10	ng	#	78

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

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