Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030116\
Data File : BM004506.D
Acq On : 02 Mar 2016 05:34
Operator : SJ/UM
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTDICCC001

Manual Integrations APPROVED
UMANGI
3/2/2016 1:19:19 PM

Quant Time: Mar 02 07:04:31 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM030116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 06:56:42 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	21273	5.00	ng		0.00
7) Naphthalene-d8	10.39	136	81466	5.00	ng		0.00
13) Acenaphthene-d10	14.26	164	41338	5.00	ng		0.00
19) Phenanthrene-d10	17.01	188	95177	5.00	ng		0.00
26) Chrysene-d12	21.23	240	73284	5.00	ng		0.00
35) Perylene-d12	23.45	264	70626	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.24	112	3931	0.81	ng		0.00
5) Phenol-d6	6.83	99	4700	0.86	ng		0.00
8) Nitrobenzene-d5	8.80	82	3415	0.84	ng		0.00
14) 2,4,6-Tribromophenol	15.79	330	1012	0.66	ng		0.00
15) 2-Fluorobiphenyl	12.88	172	11904	0.94	ng		0.00
29) Terphenyl-d14	19.67	244	11486	1.17	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.10	88	1901	0.78	ng		100
3) n-Nitrosodimethylamine	3.43	42	1355	0.65	ng		98
6) bis(2-Chloroethyl)ether	7.04	93	4215	0.74	ng		99
9) Nitrobenzene	8.84	77	3723	0.69	ng		100
10) Naphthalene	10.44	128	17516	0.70	ng		100
11) Hexachlorobutadiene	10.72	225	3051	0.83	ng		100
12) 2-Methylnaphthalene	12.08	142	10621	0.77	ng		100
16) Acenaphthylene	13.97	152	16849	0.73	ng		100
17) Acenaphthene	14.33	154	10768	0.71	ng		100
18) Fluorene	15.33	166	13541	0.74	ng		100
20) 4-Bromophenyl-phenylether	16.22	248	3035	0.60	ng	#	100
21) Hexachlorobenzene	16.35	284	3641	0.63	ng	#	100
22) Pentachlorophenol	16.73	266	409m	0.30	ng		
23) Phenanthrene	17.06	178	20183	0.62	ng		100
24) Anthracene	17.17	178	16738	0.62	ng		100
25) Fluoranthene	19.11	202	23308	0.66	ng		100
27) Benzidine	19.33	184	5481	0.71	ng		100
28) Pyrene	19.47	202	24500	0.85	ng		100
30) Benzo(a)anthracene	21.21	228	20401	0.77	ng		100
31) 3,3'-Dichlorobenzidine	21.17	252	6594	0.42	ng		100
32) Chrysene	21.27	228	23454	0.43	ng		100
33) Bis(2-ethylhexyl)phthalate	21.12	149	12019	0.61	ng	#	100
34) Indeno(1,2,3-cd)pyrene	25.71	276	20101	0.81	ng		100
36) Benzo(b)fluoranthene	22.80	252	20387	0.76	ng		100
37) Benzo(k)fluoranthene	22.84	252	20083	0.76	ng		100
38) Benzo(a)pyrene	23.36	252	18837	0.78	ng		100
39) Dibenzo(a,h)anthracene	25.71	278	15559	0.74	ng		100
40) Benzo(g,h,i)perylene	26.39	276	17760	0.77	ng		100

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

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