Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080916\
Data File : BE092236.D
Acq On : 9 Aug 2016 20:02
Operator : UM/SJ
Sample : SSTDICCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICCC0.4

Manual Integrations APPROVED
sohil
8/10/2016 6:47:33 PM

Quant Time: Aug 10 04:48:08 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 04:43:27 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	9273	5.00	ng		0.00
7) Naphthalene-d8	10.97	136	44633	5.00	ng		0.00
12) Acenaphthene-d10	14.83	164	25958	5.00	ng		0.00
18) Phenanthrene-d10	17.56	188	63185	5.00	ng		0.00
24) Chrysene-d12	21.75	240	56475	5.00	ng		0.00
31) Perylene-d12	24.13	264	58733	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.71	112	738	0.36	ng		0.00
5) Phenol-d6	7.36	99	914	0.31	ng		0.00
8) Nitrobenzene-d5	9.36	82	1064	0.34	ng		0.00
13) 2,4,6-Tribromophenol	16.33	330	240	0.31	ng		0.00
14) 2-Fluorobiphenyl	13.47	172	3112	0.39	ng		0.00
26) Terphenyl-d14	20.19	244	3534	0.37	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.44	88	507	0.43	ng		100
3) n-Nitrosodimethylamine	3.81	42	513	0.36	ng		100
6) bis(2-Chloroethyl)ether	7.59	93	853	0.36	ng		100
9) Naphthalene	11.03	128	3987	0.40	ng		100
10) Hexachlorobutadiene	11.32	225	574	0.40	ng		100
11) 2-Methylnaphthalene	12.66	142	2451	0.39	ng		100
15) Acenaphthylene	14.54	152	17210	0.38	ng		100
16) Acenaphthene	14.90	154	2833	0.39	ng		100
17) Fluorene	15.88	166	3402	0.39	ng		100
19) Hexachlorobenzene	16.89	284	1002	0.40	ng	#	100
20) Pentachlorophenol	17.24	266	173	0.26	ng		100
21) Phenanthrene	17.60	178	6296	0.39	ng		100
22) Anthracene	17.70	178	5463	0.37	ng		100
23) Fluoranthene	19.63	202	24615	0.38	ng		100
25) Pyrene	19.99	202	25528	0.38	ng		100
27) Benzo(a)anthracene	21.73	228	25260	0.37	ng		100
28) Chrysene	21.77	228	25466m	0.41	ng		
29) Bis(2-ethylhexyl)phthalate	21.63	149	2838	0.33	ng	#	100
30) Indeno(1,2,3-cd)pyrene	26.65	276	24484	0.36	ng		100
32) Benzo(b)fluoranthene	23.39	252	5682	0.37	ng		100
33) Benzo(k)fluoranthene	23.44	252	6256	0.39	ng		100
34) Benzo(a)pyrene	24.02	252	5324	0.37	ng		100
35) Dibenzo(a,h)anthracene	26.67	278	4846	0.36	ng		100
36) Benzo(g,h,i)perylene	27.42	276	22118	0.38	ng		100

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080916\
Data File : BE092236.D
Acq On : 9 Aug 2016 20:02
Operator : UM/SJ
Sample : SSTDICCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 10 04:48:08 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 04:43:27 2016
Response via : Initial Calibration

Instrument :
BNA_E
ClientSampleId :
SSTDICCC0.4

Manual Integrations APPROVED
sohil
8/10/2016 6:47:33 PM