Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE033016\
Data File : BE091558.D
Acq On : 30 Mar 2016 16:03
Operator : UM/SJ
Sample : SSTDICC0.8
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Mar 30 17:25:07 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE033016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 30 17:02:15 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	37549	5.00	ng		0.00
6) Naphthalene-d8	10.59	136	173656	5.00	ng		0.00
10) Acenaphthene-d10	14.42	164	105736	5.00	ng		0.00
16) Phenanthrene-d10	17.16	188	263449	5.00	ng		0.00
22) Chrysene-d12	21.31	240	381183	5.00	ng		0.00
28) Perylene-d12	23.76	264	325306	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.41	112	8037	1.10	ng		0.00
5) Phenol-d6	6.97	99	10704	1.07	ng		0.00
7) Nitrobenzene-d5	8.96	82	9465	1.25	ng		0.00
11) 2,4,6-Tribromophenol	15.91	330	4139	2.30	ng		0.00
12) 2-Fluorobiphenyl	13.05	172	25176	0.85	ng		0.00
24) Terphenyl-d14	19.76	244	38572	0.62	ng		-0.02
Target Compounds							Qvalue
2) 1,4-Dioxane	3.35	88	4083	0.90	ng		97
3) n-Nitrosodimethylamine	3.64	42	5679	1.11	ng		94
8) Naphthalene	10.63	128	29923	0.86	ng		97
9) 2-Methylnaphthalene	12.25	142	19448	0.89	ng		98
13) Acenaphthylene	14.14	152	36823	0.99	ng		97
14) Acenaphthene	14.49	154	22090	0.84	ng		99
15) Fluorene	15.47	166	27888	0.87	ng		100
17) Hexachlorobenzene	16.46	284	8003	0.87	ng		99
18) Pentachlorophenol	16.81	266	4320	1.42	ng	#	85
19) Phenanthrene	17.19	178	51413	0.85	ng		99
20) Anthracene	17.29	178	47113	0.94	ng		100
21) Fluoranthene	19.21	202	60372	1.07	ng		99
23) Pyrene	19.57	202	62129	0.69	ng		100
25) Benzo(a)anthracene	21.29	228	73589m	0.84	ng		
26) Chrysene	21.34	228	62182m	0.54	ng		
27) Indeno(1,2,3-cd)pyrene	26.33	276	69458	0.77	ng		99
29) Benzo(b)fluoranthene	23.01	252	65579	0.91	ng		97
30) Benzo(k)fluoranthene	23.06	252	62718	0.87	ng		95
31) Benzo(a)pyrene	23.65	252	58932	0.98	ng		95
32) Dibenzo(a,h)anthracene	26.35	278	58511	0.91	ng		98
33) Benzo(g,h,i)perylene	27.12	276	59471	0.89	ng		98

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

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