Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061818\
Data File : BE096800.D
Acq On : 18 Jun 2018 9:36
Operator : SJ/JU
Sample : SSTDICC0.1
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.1

Quant Time: Jun 18 11:40:18 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 18 11:37:15 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2550	0.40	ng		0.00
7) Naphthalene-d8	10.16	136	12395	0.40	ng		0.00
13) Acenaphthene-d10	14.03	164	6937	0.40	ng		0.00
19) Phenanthrene-d10	16.78	188	19697	0.40	ng		0.00
27) Chrysene-d12	20.98	240	13972	0.40	ng		0.00
34) Perylene-d12	23.22	264	12853	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.05	112	569	0.08	ng		0.00
5) Phenol-d6	6.60	99	805	0.08	ng		0.00
8) Nitrobenzene-d5	8.53	82	695	0.08	ng		0.00
11) 2-Methylnaphthalene-d10	11.74	152	1983	0.10	ng		0.00
14) 2,4,6-Tribromophenol	15.53	330	132	0.06	ng		0.01
15) 2-Fluorobiphenyl	12.64	172	2933	0.11	ng		0.00
25) Fluoranthene-d10	18.82	212	17224	0.09	ng		0.00
29) Terphenyl-d14	19.44	244	3327	0.11	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.10	88	446	0.108	ng	#	89
3) n-Nitrosodimethylamine	3.40	42	361	0.078	ng	#	91
6) bis(2-Chloroethyl)ether	6.85	93	959	0.095	ng		99
9) Naphthalene	10.20	128	12432	0.102	ng		99
10) Hexachlorobutadiene	10.50	225	516	0.103	ng		99
12) 2-Methylnaphthalene	11.82	142	2016	0.096	ng		98
16) Acenaphthylene	13.74	152	3125	0.093	ng		99
17) Acenaphthene	14.08	154	2190	0.099	ng		93
18) Fluorene	15.07	166	2528	0.095	ng	#	79
20) 4-Bromophenyl-phenylether	15.98	248	885	0.091	ng	#	81
21) Hexachlorobenzene	16.09	284	1136	0.103	ng	#	81
22) Pentachlorophenol	16.44	266	175	0.068	ng	#	83
23) Phenanthrene	16.81	178	5359	0.100	ng		97
24) Anthracene	16.90	178	3983	0.087	ng		97
26) Fluoranthene	18.84	202	4872	0.083	ng		98
28) Pyrene	19.21	202	4925	0.110	ng		98
30) Benzo(a)anthracene	20.97	228	3198	0.087	ng		95
31) Chrysene	21.02	228	4468	0.105	ng		98
32) Bis(2-ethylhexyl)phthalate	20.93	149	2854	0.081	ng	#	99
33) Indeno(1,2,3-cd)pyrene	25.51	276	3045	0.083	ng		93
35) Benzo(b)fluoranthene	22.54	252	3181	0.087	ng	#	95
36) Benzo(k)fluoranthene	22.58	252	3698	0.089	ng	#	90
37) Benzo(a)pyrene	23.11	252	2700	0.082	ng	#	86
38) Dibenzo(a,h)anthracene	25.53	278	2242	0.070	ng	#	91
39) Benzo(g,h,i)perylene	26.20	276	2918	0.086	ng	#	93

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

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