Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062018\
Data File : BE096833.D
Acq On : 20 Jun 2018 14:22
Operator : SJ/JU
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Jun 20 15:01:23 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 20 14:13:32 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2745	0.40	ng		0.00
7) Naphthalene-d8	10.16	136	13782	0.40	ng		0.00
13) Acenaphthene-d10	14.03	164	7785	0.40	ng		0.00
19) Phenanthrene-d10	16.78	188	21316	0.40	ng		0.00
27) Chrysene-d12	20.98	240	17914	0.40	ng		0.00
34) Perylene-d12	23.22	264	13822	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.05	112	4558	0.69	ng		0.00
5) Phenol-d6	6.59	99	6872	0.69	ng		0.00
8) Nitrobenzene-d5	8.53	82	6327	0.70	ng		0.00
11) 2-Methylnaphthalene-d10	11.73	152	14786	0.69	ng		0.00
14) 2,4,6-Tribromophenol	15.52	330	1348	0.67	ng		0.00
15) 2-Fluorobiphenyl	12.64	172	21203	0.69	ng		0.00
25) Fluoranthene-d10	18.82	212	135472	0.66	ng		0.00
29) Terphenyl-d14	19.44	244	26048	0.70	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.09	88	3206	0.724	ng	#	88
3) n-Nitrosodimethylamine	3.38	42	3623	0.768	ng	#	86
6) bis(2-Chloroethyl)ether	6.84	93	7571	0.702	ng		97
9) Naphthalene	10.20	128	92308	0.685	ng		99
10) Hexachlorobutadiene	10.50	225	3839	0.686	ng		96
12) 2-Methylnaphthalene	11.81	142	15539	0.694	ng		100
16) Acenaphthylene	13.74	152	25135	0.688	ng		100
17) Acenaphthene	14.08	154	16617	0.686	ng		95
18) Fluorene	15.07	166	19309	0.677	ng	#	79
20) 4-Bromophenyl-phenylether	15.98	248	7079	0.680	ng	#	85
21) Hexachlorobenzene	16.09	284	7753	0.650	ng	#	79
22) Pentachlorophenol	16.43	266	1431	0.660	ng	#	71
23) Phenanthrene	16.81	178	37962	0.667	ng		98
24) Anthracene	16.90	178	32336	0.687	ng		96
26) Fluoranthene	18.84	202	40755	0.670	ng		99
28) Pyrene	19.21	202	38473	0.687	ng		99
30) Benzo(a)anthracene	20.97	228	30224	0.702	ng		96
31) Chrysene	21.02	228	37090	0.700	ng		100
32) Bis(2-ethylhexyl)phthalate	20.93	149	24430	0.603	ng		99
33) Indeno(1,2,3-cd)pyrene	25.51	276	27469	0.696	ng		96
35) Benzo(b)fluoranthene	22.54	252	28422	0.686	ng	#	88
36) Benzo(k)fluoranthene	22.59	252	32782	0.700	ng		96
37) Benzo(a)pyrene	23.12	252	23456	0.675	ng	#	93
38) Dibenzo(a,h)anthracene	25.54	278	23549	0.697	ng		95
39) Benzo(g,h,i)perylene	26.21	276	25199	0.686	ng		92

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

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