Quantitation Report (Not Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070117\
Data File : BE093377.D
Acq On : 30 Jun 2017 14:36
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: Jun 30 19:50:07 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 30 19:47:49 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	4038	0.40	ng		0.00
7) Naphthalene-d8	10.74	136	20519	0.40	ng		0.00
13) Acenaphthene-d10	14.55	164	13118	0.40	ng		0.00
19) Phenanthrene-d10	17.25	188	24060	0.40	ng		-0.01
27) Chrysene-d12	21.44	240	55327	0.40	ng		0.00
34) Perylene-d12	23.95	264	50188	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.52	112	4477	0.36	ng		0.00
5) Phenol-d6	7.10	99	6612	0.35	ng		0.00
8) Nitrobenzene-d5	9.09	82	4493	0.37	ng		0.00
11) 2-Methylnaphthalene-d10	12.32	152	11787	0.36	ng		0.00
14) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng		
15) 2-Fluorobiphenyl	13.18	172	18457	0.36	ng		0.00
25) Fluoranthene-d10	19.30	212	178490	0.80	ng		0.00
29) Terphenyl-d14	19.89	244	37215	0.36	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.45	88	3054	0.366	ng		97
3) n-Nitrosodimethylamine	3.76	42	1868	0.422	ng	#	97
6) bis(2-Chloroethyl)ether	7.36	93	4777	0.361	ng	#	98
9) Naphthalene	10.79	128	75977	0.356	ng	#	73
10) Hexachlorobutadiene	11.08	225	3135	0.365	ng		99
12) 2-Methylnaphthalene	12.39	142	13031	0.359	ng		97
16) Acenaphthylene	14.27	152	20272	0.356	ng	#	87
17) Acenaphthene	14.61	154	14220	0.354	ng		99
18) Fluorene	15.59	166	19826	0.359	ng	#	87
20) 4-Bromophenyl-phenylether	16.47	248	8065	1.264	ng	#	1
21) Hexachlorobenzene	16.60	284	5390	0.586	ng	#	100
22) Pentachlorophenol	16.93	266	1615	1.387	ng	#	79
23) Phenanthrene	17.32	178	25691	0.441	ng		96
24) Anthracene	17.32	178	25682	0.438	ng	#	80
26) Fluoranthene	19.33	202	50164	0.795	ng		99
28) Pyrene	19.69	202	54658	0.374	ng	#	99
30) Benzo(a)anthracene	21.42	228	54843	0.364	ng	#	93
31) Chrysene	21.47	228	54787	0.348	ng		93
32) Bis(2-ethylhexyl)phthalate	21.35	149	86221	0.387	ng	#	67
33) Indeno(1,2,3-cd)pyrene	26.60	276	55701	0.362	ng	#	90
35) Benzo(b)fluoranthene	23.18	252	52668	0.334	ng	#	94
36) Benzo(k)fluoranthene	23.23	252	56017	0.354	ng	#	94
37) Benzo(a)pyrene	23.84	252	47757	0.341	ng	#	93
38) Dibenzo(a,h)anthracene	26.64	278	48821	0.342	ng	#	87
39) Benzo(g,h,i)perylene	27.42	276	49737	0.342	ng	#	88

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

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