Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE020419\
Data File : BE098715.D
Acq On : 4 Feb 2019 19:54
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4EC

Quant Time: Feb 04 23:59:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE012419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 24 14:44:44 2019
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.806	152	996	0.40	ng	0.00
7) Naphthalene-d8	10.602	136	4412	0.40	ng	0.00
13) Acenaphthene-d10	14.472	164	2955	0.40	ng	0.00
19) Phenanthrene-d10	17.228	188	7161	0.40	ng	0.00
27) Chrysene-d12	21.403	240	7902	0.40	ng	#-0.01
34) Perylene-d12	23.923	264	7634	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	1153	0.39	ng	0.00
5) Phenol-d6	6.989	99	1672	0.39	ng	-0.01
8) Nitrobenzene-d5	8.978	82	1652	0.38	ng	0.00
11) 2-Methylnaphthalene-d10	12.210	152	3282	0.39	ng	0.00
14) 2,4,6-Tribromophenol	15.969	330	592	0.45	ng	-0.01
15) 2-Fluorobiphenyl	13.088	172	4976	0.39	ng	-0.01
25) Fluoranthene-d10	19.253	212	37857	0.37	ng	0.00
29) Terphenyl-d14	19.856	244	8369	0.42	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.303	88	737	0.40	ng	88
3) n-Nitrosodimethylamine	3.649	42	965	0.41	ng	# 100
6) bis(2-Chloroethyl)ether	7.231	93	1139	0.39	ng	93
9) Naphthalene	10.654	128	20313	0.40	ng	100
10) Hexachlorobutadiene	10.927	225	1324	0.39	ng	96
12) 2-Methylnaphthalene	12.282	142	3299	0.39	ng	96
16) Acenaphthylene	14.198	152	5854	0.39	ng	99
17) Acenaphthene	14.529	154	3519	0.39	ng	99
18) Fluorene	15.514	166	4882	0.40	ng	95
20) 4-Bromophenyl-phenylether	16.411	248	1598	0.38	ng	95
21) Hexachlorobenzene	16.531	284	1947	0.39	ng	96
22) Pentachlorophenol	16.893	266	596	0.47	ng	93
23) Phenanthrene	17.268	178	7917	0.40	ng	99
24) Anthracene	17.361	178	6387	0.39	ng	97
26) Fluoranthene	19.288	202	10095	0.40	ng	100
28) Pyrene	19.650	202	10252	0.41	ng	99
30) Benzo(a)anthracene	21.390	228	10129	0.42	ng	98
31) Chrysene	21.451	228	10533	0.41	ng	99
32) Bis(2-ethylhexyl)phtha...	21.318	149	18147	0.41	ng	100
33) Indeno(1,2,3-cd)pyrene	26.569	276	11296	0.42	ng	99
35) Benzo(b)fluoranthene	23.150	252	9959	0.40	ng	99
36) Benzo(k)fluoranthene	23.200	252	10901	0.41	ng	99
37) Benzo(a)pyrene	23.809	252	9565	0.41	ng	98
38) Dibenzo(a,h)anthracene	26.584	278	8985	0.41	ng	95
39) Benzo(g,h,i)perylene	27.372	276	9325	0.39	ng	98

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

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