Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100841.D
Acq On : 11 Dec 2019 10:20
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: Dec 11 11:44:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	6078	0.40	ng		-0.01
7) Naphthalene-d8	10.54	136	30137	0.40	ng		0.00
13) Acenaphthene-d10	14.40	164	18801	0.40	ng		0.00
19) Phenanthrene-d10	17.12	188	68503	0.40	ng		-0.01
27) Chrysene-d12	21.31	240	67121	0.40	ng		0.00
34) Perylene-d12	23.74	264	83747	0.40	ng		-0.01
System Monitoring Compounds							
4) 2-Fluorophenol	5.36	112	6542	0.45	ng		-0.01
5) Phenol-d6	6.93	99	10442	0.48	ng		0.00
8) Nitrobenzene-d5	8.90	82	7407	0.52	ng		0.00
11) 2-Methylnaphthalene-d10	12.14	152	17117	0.39	ng		0.00
14) 2,4,6-Tribromophenol	15.88	330	2362	1.13	ng		0.00
15) 2-Fluorobiphenyl	13.02	172	24949	0.33	ng		-0.01
25) Fluoranthene-d10	19.16	212	289365	0.36	ng		0.00
29) Terphenyl-d14	19.76	244	63071	0.39	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.34	88	3900	0.352	ng		98
3) n-Nitrosodimethylamine	3.65	42	3518	0.390	ng		94
6) bis(2-Chloroethyl)ether	7.18	93	8336	0.396	ng		100
9) Naphthalene	10.59	128	122332	0.382	ng		100
10) Hexachlorobutadiene	10.89	225	4702	0.377	ng		97
12) 2-Methylnaphthalene	12.21	142	19754	0.395	ng		96
16) Acenaphthylene	14.11	152	29274	0.376	ng		100
17) Acenaphthene	14.46	154	20858	0.389	ng		98
18) Fluorene	15.42	166	33631	0.487	ng		98
20) 4-Bromophenyl-phenylether	16.32	248	11756	0.350	ng	#	68
21) Hexachlorobenzene	16.44	284	13279	0.364	ng		98
22) Pentachlorophenol	16.79	266	1694	0.671	ng	#	83
23) Phenanthrene	17.16	178	78676	0.379	ng		100
24) Anthracene	17.25	178	66309	0.382	ng		99
26) Fluoranthene	19.18	202	90004	0.363	ng		100
28) Pyrene	19.55	202	88659	0.407	ng		100
30) Benzo(a)anthracene	21.28	228	83331	0.421	ng		97
31) Chrysene	21.34	228	87517	0.376	ng		97
32) Bis(2-ethylhexyl)phthalate	21.23	149	92542	0.655	ng		100
33) Indeno(1,2,3-cd)pyrene	26.32	276	92445	0.456	ng		99
35) Benzo(b)fluoranthene	23.00	252	79640	0.347	ng		98
36) Benzo(k)fluoranthene	23.05	252	92150	0.358	ng		99
37) Benzo(a)pyrene	23.63	252	70184	0.354	ng		98
38) Dibenzo(a,h)anthracene	26.34	278	75112	0.380	ng		99
39) Benzo(g,h,i)perylene	27.10	276	78021	0.369	ng		99

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

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