Quantitation Report (Not Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
Data File : BE097905.D
Acq On : 16 Oct 2018 11:16
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: Oct 16 15:40:40 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 15:36:44 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2110	0.40	ng		0.00
7) Naphthalene-d8	10.69	136	9348	0.40	ng		0.00
13) Acenaphthene-d10	14.56	164	5967	0.40	ng		0.00
19) Phenanthrene-d10	17.30	188	15570	0.40	ng		0.00
27) Chrysene-d12	21.49	240	19787	0.40	ng		0.00
34) Perylene-d12	24.04	264	19990	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.47	112	2573	0.39	ng		0.00
5) Phenol-d6	7.06	99	3921	0.40	ng		0.00
8) Nitrobenzene-d5	9.04	82	3327	0.38	ng		0.00
11) 2-Methylnaphthalene-d10	12.30	152	6104	0.39	ng		0.00
14) 2,4,6-Tribromophenol	16.05	330	1168	0.36	ng		0.00
15) 2-Fluorobiphenyl	13.17	172	9221	0.38	ng		0.00
25) Fluoranthene-d10	19.33	212	79174	0.38	ng		0.00
29) Terphenyl-d14	19.93	244	30024	0.66	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.38	88	1298	0.353	ng		95
3) n-Nitrosodimethylamine	3.69	42	1487	0.362	ng	#	91
6) bis(2-Chloroethyl)ether	7.30	93	3143	0.376	ng		98
9) Naphthalene	10.74	128	36673	0.394	ng		100
10) Hexachlorobutadiene	11.03	225	1965	0.392	ng		97
12) 2-Methylnaphthalene	12.37	142	6235	0.391	ng		97
16) Acenaphthylene	14.27	152	10489	0.380	ng		99
17) Acenaphthene	14.61	154	6575	0.390	ng		98
18) Fluorene	15.59	166	8684	0.375	ng		100
20) 4-Bromophenyl-phenylether	16.49	248	3269	0.384	ng		97
21) Hexachlorobenzene	16.60	284	3238	0.387	ng		99
22) Pentachlorophenol	16.97	266	1010	0.483	ng		96
23) Phenanthrene	17.34	178	14674	0.371	ng		99
24) Anthracene	17.43	178	14136	0.378	ng		100
26) Fluoranthene	19.36	202	19411	0.374	ng		98
28) Pyrene	19.72	202	20987	0.368	ng		100
30) Benzo(a)anthracene	21.46	228	23129	0.381	ng		99
31) Chrysene	21.52	228	21969	0.382	ng		99
32) Bis(2-ethylhexyl)phthalate	21.39	149	66808	0.502	ng		99
33) Indeno(1,2,3-cd)pyrene	26.75	276	25632	0.406	ng		100
35) Benzo(b)fluoranthene	23.26	252	21991	0.382	ng		98
36) Benzo(k)fluoranthene	23.26	252	21991	0.367	ng		97
37) Benzo(a)pyrene	23.93	252	21881	0.390	ng		96
38) Dibenzo(a,h)anthracene	26.77	278	21489	0.391	ng		99
39) Benzo(g,h,i)perylene	27.58	276	21514	0.382	ng		100

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

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