Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE040219\
Data File : BE099273.D
Acq On : 2 Apr 2019 13:51
Operator : JU/SJ
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Apr 02 15:48:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE040219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 02 14:00:50 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	3159	0.40	ng		0.00
7) Naphthalene-d8	10.61	136	11888	0.40	ng		0.00
13) Acenaphthene-d10	14.46	164	8089	0.40	ng		0.00
19) Phenanthrene-d10	17.19	188	24251	0.40	ng		0.00
27) Chrysene-d12	21.37	240	35255	0.40	ng		0.00
34) Perylene-d12	23.86	264	36875	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.38	112	6958	0.74	ng		0.00
5) Phenol-d6	6.97	99	9552	0.66	ng		0.00
8) Nitrobenzene-d5	8.96	82	6472	0.67	ng		0.00
11) 2-Methylnaphthalene-d10	12.20	152	17298	0.76	ng		-0.01
14) 2,4,6-Tribromophenol	15.94	330	2945	1.03	ng		0.00
15) 2-Fluorobiphenyl	13.09	172	26732	0.81	ng		0.00
25) Fluoranthene-d10	19.22	212	258205	0.88	ng		0.00
29) Terphenyl-d14	19.82	244	53166	0.69	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.32	88	3532	0.848	ng		90
3) n-Nitrosodimethylamine	3.65	42	1920	0.824	ng	#	90
6) bis(2-Chloroethyl)ether	7.24	93	8518	0.752	ng		99
9) Naphthalene	10.65	128	105833	0.755	ng		100
10) Hexachlorobutadiene	10.94	225	6849	0.917	ng		99
12) 2-Methylnaphthalene	12.28	142	18192	0.726	ng		97
16) Acenaphthylene	14.18	152	31412	0.762	ng		98
17) Acenaphthene	14.53	154	18936	0.768	ng		99
18) Fluorene	15.49	166	27305	0.787	ng		98
20) 4-Bromophenyl-phenylether	16.39	248	11084	0.708	ng		97
21) Hexachlorobenzene	16.49	284	10752	0.711	ng		99
22) Pentachlorophenol	16.84	266	3906	1.960	ng		95
23) Phenanthrene	17.23	178	51697	0.726	ng		100
24) Anthracene	17.32	178	48048	0.747	ng		99
26) Fluoranthene	19.25	202	74824	0.823	ng		99
28) Pyrene	19.61	202	76850	0.656	ng		100
30) Benzo(a)anthracene	21.34	228	90947	0.782	ng		100
31) Chrysene	21.40	228	89345	0.719	ng		100
32) Bis(2-ethylhexyl)phthalate	21.27	149	108805	0.784	ng		100
33) Indeno(1,2,3-cd)pyrene	26.49	276	103797	0.882	ng		100
35) Benzo(b)fluoranthene	23.09	252	87756	0.723	ng	#	97
36) Benzo(k)fluoranthene	23.14	252	88421	0.689	ng		99
37) Benzo(a)pyrene	23.75	252	83405	0.736	ng		98
38) Dibenzo(a,h)anthracene	26.53	278	84957	0.844	ng		98
39) Benzo(g,h,i)perylene	27.31	276	84599	0.838	ng		99

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

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