Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050719\
Data File : BE099560.D
Acq On : 7 May 2019 10:49
Operator : JU/SJ
Sample : SSTDICCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICCC0.4

Quant Time: May 07 13:44:45 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 07 13:34:41 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1819	0.40	ng		0.00
7) Naphthalene-d8	10.63	136	6630	0.40	ng		0.00
13) Acenaphthene-d10	14.50	164	4894	0.40	ng		0.00
19) Phenanthrene-d10	17.24	188	13938	0.40	ng		0.00
27) Chrysene-d12	21.43	240	18601	0.40	ng		0.00
34) Perylene-d12	23.98	264	21688	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.39	112	1810	0.36	ng		0.00
5) Phenol-d6	6.98	99	2424	0.36	ng		0.00
8) Nitrobenzene-d5	8.98	82	2284	0.37	ng		0.00
11) 2-Methylnaphthalene-d10	12.22	152	4127	0.35	ng		0.00
14) 2,4,6-Tribromophenol	15.98	330	1120	0.36	ng		0.00
15) 2-Fluorobiphenyl	13.12	172	6475	0.35	ng		0.00
25) Fluoranthene-d10	19.28	212	66606	0.35	ng		0.00
29) Terphenyl-d14	19.87	244	13564	0.40	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.32	88	844	0.337	ng		100
3) n-Nitrosodimethylamine	3.64	42	614	0.407	ng	#	98
6) bis(2-Chloroethyl)ether	7.24	93	1972	0.342	ng		100
9) Naphthalene	10.68	128	25381	0.351	ng		100
10) Hexachlorobutadiene	10.95	225	1838	0.366	ng		100
12) 2-Methylnaphthalene	12.31	142	4348	0.346	ng		100
16) Acenaphthylene	14.21	152	8334	0.348	ng		100
17) Acenaphthene	14.56	154	4715	0.345	ng		100
18) Fluorene	15.53	166	6768	0.335	ng		100
20) 4-Bromophenyl-phenylether	16.42	248	2784	0.343	ng		100
21) Hexachlorobenzene	16.55	284	2848	0.368	ng		100
22) Pentachlorophenol	16.89	266	1043	0.299	ng		100
23) Phenanthrene	17.28	178	12383	0.344	ng		100
24) Anthracene	17.38	178	11600	0.338	ng		100
26) Fluoranthene	19.31	202	18973	0.352	ng		100
28) Pyrene	19.67	202	19685	0.406	ng		100
30) Benzo(a)anthracene	21.41	228	22642	0.382	ng		100
31) Chrysene	21.46	228	22711	0.392	ng		100
32) Bis(2-ethylhexyl)phthalate	21.33	149	35594	0.345	ng		100
33) Indeno(1,2,3-cd)pyrene	26.68	276	26627	0.388	ng		100
35) Benzo(b)fluoranthene	23.19	252	21846	0.349	ng		100
36) Benzo(k)fluoranthene	23.24	252	21353	0.354	ng		100
37) Benzo(a)pyrene	23.86	252	20948	0.354	ng		100
38) Dibenzo(a,h)anthracene	26.71	278	21621	0.358	ng		100
39) Benzo(g,h,i)perylene	27.51	276	21969	0.363	ng		100

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

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