Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032519\
Data File : BE099181.D
Acq On : 25 Mar 2019 13:46
Operator : JU/SJ
Sample : SSTDICC3.2
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC3.2

Quant Time: Mar 25 14:26:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 25 14:07:14 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	3945	0.40	ng		-0.01
7) Naphthalene-d8	10.63	136	17778	0.40	ng		0.00
13) Acenaphthene-d10	14.47	164	13190	0.40	ng		-0.01
19) Phenanthrene-d10	17.21	188	33209	0.40	ng		0.00
27) Chrysene-d12	21.38	240	40091	0.40	ng		-0.01
34) Perylene-d12	23.89	264	38923	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.41	112	36557	3.11	ng		0.00
5) Phenol-d6	6.99	99	57710	3.41	ng		0.00
8) Nitrobenzene-d5	8.99	82	49907	6.42	ng		0.00
11) 2-Methylnaphthalene-d10	12.22	152	109485	3.42	ng		0.00
14) 2,4,6-Tribromophenol	15.96	330	21849	5.63	ng		0.00
15) 2-Fluorobiphenyl	13.10	172	167810	3.20	ng		0.00
25) Fluoranthene-d10	19.24	212	1341017	3.03	ng		0.00
29) Terphenyl-d14	19.83	244	279602	3.30	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.34	88	16024	2.618	ng		96
3) n-Nitrosodimethylamine	3.65	42	10826	3.428	ng	#	90
6) bis(2-Chloroethyl)ether	7.25	93	45836	3.335	ng		98
9) Naphthalene	10.68	128	664628	3.172	ng		100
10) Hexachlorobutadiene	10.95	225	35781	3.493	ng		96
12) 2-Methylnaphthalene	12.30	142	121378	3.383	ng		98
16) Acenaphthylene	14.20	152	228868	3.251	ng		100
17) Acenaphthene	14.54	154	130417	3.124	ng		98
18) Fluorene	15.51	166	182397	2.995	ng		99
20) 4-Bromophenyl-phenylether	16.41	248	71383	3.883	ng		98
21) Hexachlorobenzene	16.52	284	66206	3.878	ng		99
22) Pentachlorophenol	16.85	266	13587	2.194	ng	#	88
23) Phenanthrene	17.25	178	309952	3.143	ng		100
24) Anthracene	17.33	178	301375	3.230	ng		100
26) Fluoranthene	19.27	202	411156	2.963	ng		99
28) Pyrene	19.63	202	416478	3.063	ng		100
30) Benzo(a)anthracene	21.37	228	428107	3.031	ng		100
31) Chrysene	21.42	228	439086	3.111	ng		99
32) Bis(2-ethylhexyl)phthalate	21.29	149	542019	2.957	ng		100
33) Indeno(1,2,3-cd)pyrene	26.55	276	447071	2.946	ng		98
35) Benzo(b)fluoranthene	23.12	252	418401	3.107	ng		99
36) Benzo(k)fluoranthene	23.17	252	440009	3.335	ng		99
37) Benzo(a)pyrene	23.78	252	394176	3.167	ng		98
38) Dibenzo(a,h)anthracene	26.58	278	372826	3.081	ng		98
39) Benzo(g,h,i)perylene	27.37	276	365294	2.939	ng		100

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

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