Quantitation Report (QT Reviewed)

Data Path : S:\HPCHEM1\BNA_E\DATA\BE050517\
Data File : BE092928.D
Acq On : 5 May 2017 13:33
Operator : SJ/MA
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE050517

Quant Time: May 05 14:08:13 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE050517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 05 14:05:07 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	738	0.40	ng		0.00
7) Naphthalene-d8	10.53	136	3028	0.40	ng		0.00
13) Acenaphthene-d10	14.39	164	1367	0.40	ng		0.00
19) Phenanthrene-d10	17.13	188	3649	0.40	ng		0.00
26) Chrysene-d12	21.28	240	3278	0.40	ng		0.00
33) Perylene-d12	23.69	264	2485	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.35	112	640	0.39	ng		0.00
5) Phenol-d6	6.94	99	878	0.38	ng		0.00
8) Nitrobenzene-d5	8.90	82	821	0.40	ng		0.00
11) 2-Methylnaphthalene-d10	12.13	152	1769	0.39	ng		0.00
14) 2,4,6-Tribromophenol	15.87	330	81	0.38	ng		0.00
15) 2-Fluorobiphenyl	13.01	172	2530	0.40	ng		0.00
24) Fluoranthene-d10	19.15	212	3193	0.38	ng		0.00
28) Terphenyl-d14	19.74	244	2387	0.42	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.31	88	544	0.38	ng		98
3) n-Nitrosodimethylamine	3.61	42	411	0.38	ng	#	84
6) bis(2-Chloroethyl)ether	7.20	93	1023	0.38	ng		98
9) Naphthalene	10.58	128	3208	0.40	ng	#	87
10) Hexachlorobutadiene	10.89	225	438	0.40	ng		97
12) 2-Methylnaphthalene	12.19	142	1922	0.39	ng		96
16) Acenaphthylene	14.10	152	8164	0.38	ng		99
17) Acenaphthene	14.44	154	1697	0.39	ng		94
18) Fluorene	15.43	166	2043	0.38	ng		99
20) Hexachlorobenzene	16.47	284	594	0.38	ng	#	100
21) Pentachlorophenol	16.79	266	106	0.42	ng	#	84
22) Phenanthrene	17.18	178	3859	0.38	ng		98
23) Anthracene	17.27	178	2900	0.37	ng		99
25) Fluoranthene	19.17	202	16531	0.37	ng	#	98
27) Pyrene	19.53	202	17074	0.40	ng	#	98
29) Benzo(a)anthracene	21.25	228	2991	0.39	ng	#	87
30) Chrysene	21.30	228	4081	0.42	ng	#	89
31) Bis(2-ethylhexyl)phthalate	21.20	149	817	0.39	ng	#	97
32) Indeno(1,2,3-cd)pyrene	26.22	276	12783	0.36	ng		99
34) Benzo(b)fluoranthene	22.95	252	3235	0.40	ng	#	84
35) Benzo(k)fluoranthene	23.00	252	3268	0.40	ng	#	80
36) Benzo(a)pyrene	23.57	252	2398	0.36	ng	#	77
37) Dibenzo(a,h)anthracene	26.25	278	2848	0.38	ng	#	78
38) Benzo(g,h,i)perylene	26.99	276	12674	0.40	ng	#	82

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

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