Quantitation Report (QT Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA E\DATA\BE090517\
Data File : BE093928.D
Acq On : 5 Sep 2017 9:29
Operator : SJ/JU
Sample : SSTDICC0.1
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDICC0.1

Manual Integrations APPROVED
Sohil
9/7/2017 6:30:38 PM

Quant Time: Sep 05 13:31:35 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 12:08:29 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1812	0.40	ng		0.00
7) Naphthalene-d8	10.72	136	9370	0.40	ng		0.02
13) Acenaphthene-d10	14.52	164	5294	0.40	ng		0.00
19) Phenanthrene-d10	17.27	188	11109	0.40	ng		0.00
27) Chrysene-d12	21.43	240	13814	0.40	ng		0.00
34) Perylene-d12	23.94	264	11499	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.53	112	588	0.11	ng		0.00
5) Phenol-d6	7.12	99	830	0.10	ng		0.01
8) Nitrobenzene-d5	9.08	82	741	0.09	ng		0.00
11) 2-Methylnaphthalene-d10	12.31	152	1477	0.10	ng		0.01
14) 2,4,6-Tribromophenol	16.03	330	190	0.13	ng		0.00
15) 2-Fluorobiphenyl	13.16	172	2213	0.11	ng		0.00
25) Fluoranthene-d10	19.29	212	15128	0.12	ng		0.00
29) Terphenyl-d14	19.87	244	2629	0.11	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.42	88	242m	0.097	ng		
3) n-Nitrosodimethylamine	3.75	42	255	0.092	ng	#	82
6) bis(2-Chloroethyl)ether	7.35	93	709	0.108	ng		78
9) Naphthalene	10.76	128	10805	0.101	ng	#	97
10) Hexachlorobutadiene	11.04	225	414	0.110	ng		94
12) 2-Methylnaphthalene	12.38	142	1651	0.093	ng		96
16) Acenaphthylene	14.26	152	2497	0.095	ng		99
17) Acenaphthene	14.60	154	1793	0.101	ng		99
18) Fluorene	15.58	166	2338	0.102	ng		100
20) 4-Bromophenyl-phenylether	16.45	248	708	0.154	ng	#	68
21) Hexachlorobenzene	16.58	284	706	0.176	ng	#	100
23) Phenanthrene	17.30	178	3933m	0.139	ng		
24) Anthracene	17.40	178	3354m	0.100	ng		
26) Fluoranthene	19.32	202	4502	0.116	ng		99
28) Pyrene	19.67	202	4723	0.102	ng		98
30) Benzo(a)anthracene	21.40	228	4096	0.100	ng	#	90
31) Chrysene	21.46	228	4546	0.100	ng	#	93
32) Bis(2-ethylhexyl)phthalate	21.30	149	11917	0.138	ng	#	99
33) Indeno(1,2,3-cd)pyrene	26.61	276	2963	0.095	ng	#	63
35) Benzo(b)fluoranthene	23.17	252	3421	0.093	ng	#	70
36) Benzo(k)fluoranthene	23.22	252	4274	0.102	ng	#	69
37) Benzo(a)pyrene	23.83	252	3638	0.100	ng	#	43
38) Dibenzo(a,h)anthracene	26.63	278	2215	0.086	ng	#	48
39) Benzo(g,h,i)perylene	27.43	276	2542	0.091	ng	#	73

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

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