Quantitation Report (QT Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA E\DATA\BE090517\
Data File : BE093931.D
Acq On : 5 Sep 2017 11:19
Operator : SJ/JU
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDICC0.8

Manual Integrations APPROVED
Sohil
9/7/2017 6:30:45 PM

Quant Time: Sep 05 13:18:51 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	1825	0.40	ng		0.00
7) Naphthalene-d8	10.71	136	9541	0.40	ng		0.00
13) Acenaphthene-d10	14.53	164	5493	0.40	ng		0.00
19) Phenanthrene-d10	17.24	188	9028	0.40	ng		-0.03
27) Chrysene-d12	21.43	240	13898	0.40	ng		0.00
34) Perylene-d12	23.94	264	11353	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.53	112	4581	0.82	ng		0.00
5) Phenol-d6	7.12	99	6876	0.80	ng		0.00
8) Nitrobenzene-d5	9.08	82	6267	0.79	ng		0.00
11) 2-Methylnaphthalene-d10	12.29	152	12163	0.81	ng		0.00
14) 2,4,6-Tribromophenol	16.03	330	1385	0.94	ng		0.00
15) 2-Fluorobiphenyl	13.16	172	16948	0.78	ng		0.00
25) Fluoranthene-d10	19.28	212	121015	1.16	ng		0.00
29) Terphenyl-d14	19.87	244	20250	0.81	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.43	88	1904	0.759	ng		94
3) n-Nitrosodimethylamine	3.74	42	2334	0.836	ng	#	98
6) bis(2-Chloroethyl)ether	7.34	93	5645	0.853	ng		92
9) Naphthalene	10.76	128	85541	0.787	ng		100
10) Hexachlorobutadiene	11.04	225	3048	0.795	ng		98
12) 2-Methylnaphthalene	12.36	142	14547	0.801	ng		98
16) Acenaphthylene	14.24	152	21700	0.798	ng		100
17) Acenaphthene	14.58	154	14649	0.799	ng		98
18) Fluorene	15.58	166	18977	0.800	ng		99
20) 4-Bromophenyl-phenylether	16.45	248	4025	1.077	ng	#	95
21) Hexachlorobenzene	16.58	284	4406	1.353	ng	#	100
22) Pentachlorophenol	16.94	266	1170	0.827	ng		97
23) Phenanthrene	17.30	178	23313m	1.016	ng		
24) Anthracene	17.40	178	22038m	0.805	ng		
26) Fluoranthene	19.31	202	35843	1.140	ng		100
28) Pyrene	19.67	202	37945	0.812	ng		99
30) Benzo(a)anthracene	21.40	228	34395	0.833	ng		99
31) Chrysene	21.46	228	35039	0.764	ng		98
32) Bis(2-ethylhexyl)phthalate	21.30	149	76103	0.875	ng		100
33) Indeno(1,2,3-cd)pyrene	26.61	276	23688	0.757	ng		99
35) Benzo(b)fluoranthene	23.16	252	28211	0.778	ng		98
36) Benzo(k)fluoranthene	23.21	252	36034	0.870	ng		99
37) Benzo(a)pyrene	23.82	252	29168	0.811	ng		99
38) Dibenzo(a,h)anthracene	26.63	278	19536	0.769	ng		99
39) Benzo(g,h,i)perylene	27.42	276	20375	0.737	ng		98

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

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