Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082617\
Data File : BE093887.D
Acq On : 26 Aug 2017 13:26
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDICV0.4

Manual Integrations APPROVED
Sohil
8/28/2017 3:12:31 PM

Quant Time: Aug 28 11:29:02 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE082617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 28 11:26:07 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1879	0.40	ng		0.00
7) Naphthalene-d8	10.71	136	9738	0.40	ng		0.00
13) Acenaphthene-d10	14.52	164	5420	0.40	ng		0.00
19) Phenanthrene-d10	17.23	188	10635	0.40	ng		0.00
27) Chrysene-d12	21.42	240	12246	0.40	ng		0.00
34) Perylene-d12	23.93	264	10111	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.52	112	2300	0.40	ng		0.00
5) Phenol-d6	7.09	99	3365	0.38	ng		0.00
8) Nitrobenzene-d5	9.06	82	3018	0.37	ng		0.00
11) 2-Methylnaphthalene-d10	12.29	152	6085	0.40	ng		0.00
14) 2,4,6-Tribromophenol	16.03	330	477	0.33	ng		0.00
15) 2-Fluorobiphenyl	13.15	172	8482	0.39	ng		-0.01
25) Fluoranthene-d10	19.28	212	51629	0.42	ng		0.00
29) Terphenyl-d14	19.87	244	8569	0.39	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.43	88	1035	0.401	ng		98
3) n-Nitrosodimethylamine	3.75	42	1114	0.388	ng		99
6) bis(2-Chloroethyl)ether	7.34	93	2767	0.406	ng		93
9) Naphthalene	10.76	128	43290	0.390	ng		100
10) Hexachlorobutadiene	11.04	225	1546	0.395	ng		97
12) 2-Methylnaphthalene	12.36	142	7187	0.388	ng		99
16) Acenaphthylene	14.24	152	10344	0.386	ng		99
17) Acenaphthene	14.58	154	6967	0.385	ng		99
18) Fluorene	15.56	166	8908	0.381	ng		100
20) 4-Bromophenyl-phenylether	16.45	248	1940	0.495	ng		92
21) Hexachlorobenzene	16.58	284	1625	0.460	ng	#	100
22) Pentachlorophenol	16.94	266	491	0.295	ng		97
23) Phenanthrene	17.30	178	10412	0.385	ng	#	99
24) Anthracene	17.36	178	11589m	0.359	ng		
26) Fluoranthene	19.31	202	15539	0.420	ng		100
28) Pyrene	19.67	202	16053	0.390	ng		99
30) Benzo(a)anthracene	21.40	228	13680	0.376	ng		99
31) Chrysene	21.45	228	15977	0.395	ng		99
32) Bis(2-ethylhexyl)phthalate	21.30	149	26248	0.342	ng		100
33) Indeno(1,2,3-cd)pyrene	26.59	276	9996	0.362	ng		99
35) Benzo(b)fluoranthene	23.15	252	11710	0.363	ng		99
36) Benzo(k)fluoranthene	23.20	252	14088	0.382	ng		98
37) Benzo(a)pyrene	23.81	252	12092	0.377	ng		99
38) Dibenzo(a,h)anthracene	26.61	278	8244	0.365	ng		99
39) Benzo(g,h,i)perylene	27.40	276	9145	0.371	ng		99

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

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