Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080117\
Data File : BE093615.D
Acq On : 1 Aug 2017 14:14
Operator : SJ/JU
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Aug 01 14:52:29 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 01 14:41:34 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2338	0.40	ng		0.00
7) Naphthalene-d8	10.71	136	10633	0.40	ng		0.00
13) Acenaphthene-d10	14.53	164	6082	0.40	ng		0.00
19) Phenanthrene-d10	17.24	188	17782	0.40	ng		0.00
27) Chrysene-d12	21.40	240	18949	0.40	ng		0.00
34) Perylene-d12	23.91	264	15166	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.51	112	5382	0.66	ng		0.00
5) Phenol-d6	7.09	99	7780	0.68	ng		0.00
8) Nitrobenzene-d5	9.06	82	6351	0.75	ng		0.00
11) 2-Methylnaphthalene-d10	12.29	152	13366	0.78	ng		0.00
14) 2,4,6-Tribromophenol	16.00	330	1134	0.51	ng		0.00
15) 2-Fluorobiphenyl	13.15	172	19139	0.80	ng		-0.01
25) Fluoranthene-d10	19.27	212	149218	0.53	ng		0.00
29) Terphenyl-d14	19.86	244	26544	0.77	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.44	88	3052	0.786	ng		97
3) n-Nitrosodimethylamine	3.74	42	2966	0.772	ng		97
6) bis(2-Chloroethyl)ether	7.34	93	6949	0.768	ng	#	99
9) Naphthalene	10.76	128	96709	0.781	ng		99
10) Hexachlorobutadiene	11.04	225	3658	0.760	ng		98
12) 2-Methylnaphthalene	12.36	142	16012	0.771	ng		97
16) Acenaphthylene	14.24	152	23048	0.752	ng	#	88
17) Acenaphthene	14.59	154	16202	0.779	ng		99
18) Fluorene	15.56	166	21606	0.738	ng	#	87
20) 4-Bromophenyl-phenylether	16.42	248	3639	0.525	ng	#	1
21) Hexachlorobenzene	16.55	284	3962	0.595	ng	#	100
22) Pentachlorophenol	16.91	266	1762	0.453	ng		95
23) Phenanthrene	17.27	178	29282	0.726	ng	#	79
24) Anthracene	17.37	178	41826	0.678	ng	#	89
26) Fluoranthene	19.30	202	46074	0.536	ng		99
28) Pyrene	19.66	202	47198	0.769	ng	#	99
30) Benzo(a)anthracene	21.39	228	44213	0.725	ng		95
31) Chrysene	21.45	228	48739	0.807	ng		96
32) Bis(2-ethylhexyl)phthalate	21.30	149	54669	0.414	ng		98
33) Indeno(1,2,3-cd)pyrene	26.55	276	43497	0.761	ng		93
35) Benzo(b)fluoranthene	23.14	252	42914	0.806	ng	#	93
36) Benzo(k)fluoranthene	23.19	252	44035	0.837	ng	#	92
37) Benzo(a)pyrene	23.80	252	38843	0.784	ng	#	91
38) Dibenzo(a,h)anthracene	26.57	278	38057	0.813	ng	#	89
39) Benzo(g,h,i)perylene	27.36	276	38196	0.813	ng	#	91

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

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