Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE111618\
Data File : BE098233.D
Acq On : 16 Nov 2018 17:16
Operator : SJ/JU
Sample : SSTDICC0.8
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Nov 16 17:47:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 16 17:10:30 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	1356	0.40	ng		0.00
7) Naphthalene-d8	10.67	136	6791	0.40	ng		0.00
13) Acenaphthene-d10	14.53	164	4345	0.40	ng		0.00
19) Phenanthrene-d10	17.28	188	9954	0.40	ng		0.00
27) Chrysene-d12	21.46	240	10763	0.40	ng		0.00
34) Perylene-d12	24.02	264	10058	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.45	112	2790	0.73	ng		-0.01
5) Phenol-d6	7.03	99	4507	0.73	ng		0.00
8) Nitrobenzene-d5	9.02	82	4934	0.73	ng		0.00
11) 2-Methylnaphthalene-d10	12.27	152	9189	0.77	ng		0.00
14) 2,4,6-Tribromophenol	16.02	330	1166	0.62	ng		-0.01
15) 2-Fluorobiphenyl	13.15	172	14162	0.78	ng		0.00
25) Fluoranthene-d10	19.31	212	94271	0.76	ng		0.00
29) Terphenyl-d14	19.91	244	19016	0.78	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.37	88	1722	0.781	ng		89
3) n-Nitrosodimethylamine	3.68	42	1970	0.740	ng	#	85
6) bis(2-Chloroethyl)ether	7.29	93	3838	0.773	ng		100
9) Naphthalene	10.72	128	53195	0.770	ng		99
10) Hexachlorobutadiene	11.00	225	3086	0.719	ng		98
12) 2-Methylnaphthalene	12.34	142	9513	0.794	ng		92
16) Acenaphthylene	14.26	152	16228	0.815	ng		99
17) Acenaphthene	14.60	154	9679	0.796	ng		99
18) Fluorene	15.57	166	12301	0.790	ng		100
20) 4-Bromophenyl-phenylether	16.47	248	4861	0.801	ng		91
21) Hexachlorobenzene	16.58	284	4878	0.773	ng		98
22) Pentachlorophenol	16.94	266	761	0.694	ng		93
23) Phenanthrene	17.32	178	19763	0.797	ng		100
24) Anthracene	17.41	178	18432	0.798	ng		98
26) Fluoranthene	19.34	202	23993	0.782	ng		98
28) Pyrene	19.70	202	24538	0.828	ng		99
30) Benzo(a)anthracene	21.45	228	24145	0.790	ng		99
31) Chrysene	21.50	228	25081	0.808	ng		98
32) Bis(2-ethylhexyl)phthalate	21.37	149	36216	0.570	ng		100
33) Indeno(1,2,3-cd)pyrene	26.70	276	22100	0.776	ng	#	91
35) Benzo(b)fluoranthene	23.23	252	22288	0.781	ng		96
36) Benzo(k)fluoranthene	23.28	252	24728	0.821	ng		100
37) Benzo(a)pyrene	23.90	252	21764	0.795	ng		98
38) Dibenzo(a,h)anthracene	26.74	278	17575	0.754	ng	#	92
39) Benzo(g,h,i)perylene	27.53	276	19147	0.787	ng		97

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

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