

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE050118\
 Data File : BE096287.D
 Acq On : 2 May 2018 9:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4EC

Quant Time: May 02 11:19:58 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 30 14:27:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	869	0.40	ng	0.00
7) Naphthalene-d8	11.19	136	3311	0.40	ng	0.00
13) Acenaphthene-d10	14.94	164	1810	0.40	ng	0.00
19) Phenanthrene-d10	17.65	188	4173	0.40	ng	0.00
27) Chrysene-d12	21.74	240	4213	0.40	ng	0.00
34) Perylene-d12	24.35	264	3927	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.87	112	942	0.35	ng	0.00
5) Phenol-d6	7.52	99	1407	0.38	ng	0.00
8) Nitrobenzene-d5	9.54	82	1476	0.38	ng	0.00
11) 2-Methylnaphthalene-d10	12.73	152	1936	0.37	ng	0.00
14) 2,4,6-Tribromophenol	16.41	330	274	0.31	ng	0.00
15) 2-Fluorobiphenyl	13.57	172	3032	0.39	ng	0.00
25) Fluoranthene-d10	19.63	212	20228	0.36	ng	0.00
29) Terphenyl-d14	20.19	244	3467	0.37	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.60	88	450	0.335	ng	92
3) n-Nitrosodimethylamine	3.97	42	647	0.393	ng	# 72
6) bis(2-Chloroethyl)ether	7.76	93	1174	0.360	ng	93
9) Naphthalene	11.24	128	12998	0.372	ng	100
10) Hexachlorobutadiene	11.50	225	491	0.317	ng	88
12) 2-Methylnaphthalene	12.81	142	2158	0.374	ng	96
16) Acenaphthylene	14.67	152	3722	0.385	ng	100
17) Acenaphthene	15.00	154	2176	0.375	ng	92
18) Fluorene	15.97	166	2785	0.387	ng	# 65
20) 4-Bromophenyl-phenylether	16.83	248	959	0.387	ng	# 77
21) Hexachlorobenzene	16.96	284	977	0.396	ng	# 80
22) Pentachlorophenol	17.31	266	366	0.457	ng	84
23) Phenanthrene	17.68	178	4483	0.382	ng	99
24) Anthracene	17.78	178	4265	0.381	ng	# 95
26) Fluoranthene	19.66	202	5635	0.366	ng	98
28) Pyrene	20.00	202	403	0.049	ng	# 86
30) Benzo(a)anthracene	21.73	228	4996	0.370	ng	97
31) Chrysene	21.79	228	5203	0.399	ng	99
32) Bis(2-ethylhexyl)phthalate	21.60	149	14078	0.404	ng	100
33) Indeno(1,2,3-cd)pyrene	27.13	276	4456	0.352	ng	94
35) Benzo(b)fluoranthene	23.54	252	4544	0.386	ng	# 94
36) Benzo(k)fluoranthene	23.59	252	4731	0.390	ng	# 90
37) Benzo(a)pyrene	24.23	252	4340	0.383	ng	# 92
38) Dibenzo(a,h)anthracene	27.14	278	3597	0.360	ng	94
39) Benzo(g,h,i)perylene	27.99	276	3917	0.371	ng	# 93

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

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