

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022618\
 Data File : BE095688.D
 Acq On : 26 Feb 2018 14:42
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Feb 26 17:49:51 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE022318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 12:12:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1324	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	5396	0.40	ng	-0.02
13) Acenaphthene-d10	15.00	164	3200	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	7427	0.40	ng	0.01
27) Chrysene-d12	21.79	240	8042	0.40	ng	0.00
34) Perylene-d12	24.42	264	7373	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.92	112	1469	0.42	ng	0.00
5) Phenol-d6	7.56	99	2031	0.42	ng	0.00
8) Nitrobenzene-d5	9.61	82	2122	0.38	ng	0.00
11) 2-Methylnaphthalene-d10	12.80	152	3630	0.40	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	419	0.39	ng	0.00
15) 2-Fluorobiphenyl	13.64	172	5494	0.38	ng	0.00
25) Fluoranthene-d10	19.69	212	37341	0.38	ng	0.00
29) Terphenyl-d14	20.25	244	6260	0.39	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.64	88	832	0.399	ng	93
3) n-Nitrosodimethylamine	3.99	42	1110	0.429	ng	80
6) bis(2-Chloroethyl)ether	7.82	93	2131	0.429	ng	95
9) Naphthalene	11.31	128	24916	0.401	ng	99
10) Hexachlorobutadiene	11.57	225	1503	0.414	ng	98
12) 2-Methylnaphthalene	12.88	142	4280	0.399	ng	98
16) Acenaphthylene	14.74	152	6956	0.387	ng	100
17) Acenaphthene	15.06	154	4305	0.381	ng	94
18) Fluorene	16.02	166	5453	0.390	ng	# 79
20) 4-Bromophenyl-phenylether	16.89	248	1719	0.369	ng	# 77
21) Hexachlorobenzene	17.03	284	1768	0.366	ng	# 81
22) Pentachlorophenol	17.36	266	403	0.406	ng	# 77
23) Phenanthrene	17.74	178	8820	0.380	ng	99
24) Anthracene	17.83	178	7947	0.373	ng	95
26) Fluoranthene	19.72	202	11098	0.389	ng	# 97
28) Pyrene	20.06	202	10833	0.329	ng	# 90
30) Benzo(a)anthracene	21.78	228	10229	0.390	ng	97
31) Chrysene	21.84	228	10017	0.384	ng	99
32) Bis(2-ethylhexyl)phthalate	21.65	149	18681	0.403	ng	100
33) Indeno(1,2,3-cd)pyrene	27.23	276	9886	0.391	ng	# 93
35) Benzo(b)fluoranthene	23.61	252	9775	0.372	ng	# 90
36) Benzo(k)fluoranthene	23.66	252	9839	0.377	ng	96
37) Benzo(a)pyrene	24.30	252	9212	0.389	ng	93
38) Dibenzo(a,h)anthracene	27.25	278	7977	0.379	ng	97
39) Benzo(g,h,i)perylene	28.10	276	8472	0.375	ng	# 92

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

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