

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032318\
 Data File : BE095815.D
 Acq On : 23 Mar 2018 8:46
 Operator : SJ/JU
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.1

Quant Time: Mar 23 10:57:46 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 10:57:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	981	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	3818	0.40	ng	0.00
13) Acenaphthene-d10	15.00	164	2309	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	5876	0.40	ng	0.00
27) Chrysene-d12	21.80	240	6988	0.40	ng	0.00
34) Perylene-d12	24.44	264	6674	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.92	112	341	0.13	ng	0.00
5) Phenol-d6	7.57	99	441	0.12	ng	0.00
8) Nitrobenzene-d5	9.61	82	453	0.12	ng	0.01
11) 2-Methylnaphthalene-d10	12.80	152	657	0.10	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	134	0.16	ng	0.00
15) 2-Fluorobiphenyl	13.64	172	967	0.09	ng	0.00
25) Fluoranthene-d10	19.69	212	8877	0.12	ng	0.00
29) Terphenyl-d14	20.24	244	1519	0.11	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.63	88	176	0.114	ng	92
3) n-Nitrosodimethylamine	3.99	42	235	0.123	ng	# 70
6) bis(2-Chloroethyl)ether	7.82	93	410	0.111	ng	# 87
9) Naphthalene	11.31	128	4491	0.102	ng	# 96
10) Hexachlorobutadiene	11.56	225	309	0.120	ng	96
12) 2-Methylnaphthalene	12.87	142	750	0.099	ng	# 92
16) Acenaphthylene	14.73	152	1308	0.101	ng	98
17) Acenaphthene	15.06	154	785	0.096	ng	93
18) Fluorene	16.03	166	1062	0.105	ng	# 80
20) 4-Bromophenyl-phenylether	16.89	248	348	0.094	ng	# 62
21) Hexachlorobenzene	17.03	284	356	0.093	ng	# 84
22) Pentachlorophenol	17.37	266	110	0.125	ng	# 81
23) Phenanthrene	17.75	178	1744	0.095	ng	97
24) Anthracene	17.83	178	1740	0.103	ng	# 94
26) Fluoranthene	19.72	202	2455	0.109	ng	95
28) Pyrene	20.07	202	2641	0.092	ng	# 83
30) Benzo(a)anthracene	21.78	228	2389	0.105	ng	# 85
31) Chrysene	21.84	228	2236	0.099	ng	# 83
32) Bis(2-ethylhexyl)phthalate	21.64	149	9208	0.228	ng	# 98
33) Indeno(1,2,3-cd)pyrene	27.25	276	2350	0.107	ng	92
35) Benzo(b)fluoranthene	23.62	252	2250	0.095	ng	# 63
36) Benzo(k)fluoranthene	23.67	252	2172	0.092	ng	# 44
37) Benzo(a)pyrene	24.31	252	2047	0.096	ng	# 42
38) Dibenzo(a,h)anthracene	27.27	278	1931	0.101	ng	# 49
39) Benzo(g,h,i)perylene	28.13	276	2009	0.098	ng	# 59

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

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