

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020515\
 Data File : BE089506.D
 Acq On : 5 Feb 2015 15:23
 Operator : TP/IZ
 Sample : SSTDICC0.2
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.2

Quant Time: Feb 06 03:15:12 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 06 02:40:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.49	152	26545	5.00	ng	0.00
2) Naphthalene-d8	12.40	136	106754	5.00	ng	0.00
6) Acenaphthene-d10	16.59	164	47143	5.00	ng	0.00
11) Phenanthrene-d10	19.91	188	89596	5.00	ng	0.00
16) Chrysene-d12	23.76	240	95581	5.00	ng	0.00
22) Perylene-d12	25.46	264	95807	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.76	82	1613	0.20	ng	0.00
7) 2-Fluorobiphenyl	15.03	172	3043	0.20	ng	0.00
18) Terphenyl-d14	22.40	244	3079	0.19	ng	0.00
Target Compounds						Ovalue
4) Naphthalene	12.44	128	5632	0.21	ng	99
5) 2-Methylnaphthalene	14.10	142	3471	0.20	ng	97
8) Acenaphthylene	16.24	152	19162	0.19	ng	99
9) Acenaphthene	16.65	154	3137	0.20	ng	97
10) Fluorene	17.91	166	3173	0.18	ng	95
13) Phenanthrene	19.94	178	5406	0.20	ng	98
14) Anthracene	20.03	178	4284	0.18	ng	97
15) Fluoranthene	21.83	202	4754	0.18	ng	100
17) Pyrene	22.15	202	5083	0.19	ng	99
19) Benzo(a)anthracene	23.75	228	19945	0.19	ng	99
20) Chrysene	23.79	228	21047	0.20	ng	98
21) Indeno(1,2,3-cd)pyrene	26.92	276	19081	0.17	ng	99
23) Benzo(b)fluoranthene	25.02	252	3679	0.17	ng	# 96
24) Benzo(k)fluoranthene	25.05	252	5989	0.20	ng	97
25) Benzo(a)pyrene	25.39	252	3809	0.17	ng	# 93
26) Dibenzo(a,h)anthracene	26.95	278	4254	0.18	ng	# 90
27) Benzo(g,h,i)perylene	27.36	276	18070	0.18	ng	94

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

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