

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041815\
 Data File : BM001041.D
 Acq On : 17 Apr 2015 22:08
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Quant Time: Apr 18 07:45:02 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM041815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 18 05:45:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	16286	5.00	ng	0.00
4) Naphthalene-d8	10.36	136	58053	5.00	ng	0.00
8) Acenaphthene-d10	14.24	164	30087	5.00	ng	0.00
13) Phenanthrene-d10	16.99	188	65552	5.00	ng	0.00
19) Chrysene-d12	21.19	240	55279	5.00	ng	0.00
25) Perylene-d12	23.36	264	50580	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.75	82	263	0.09	ng	0.01
9) 2-Fluorobiphenyl	12.86	172	980	0.10	ng	0.00
21) Terphenyl-d14	19.63	244	723	0.11	ng	0.00
Target Compounds						Ovalue
6) Naphthalene	10.41	128	1396	0.10	ng	# 88
7) 2-Methylnaphthalene	12.05	142	858	0.10	ng	# 94
10) Acenaphthylene	13.95	152	1192	0.09	ng	99
11) Acenaphthene	14.29	154	1018	0.11	ng	91
12) Fluorene	15.30	166	1044	0.10	ng	100
14) Hexachlorobenzene	16.31	284	361	0.10	ng	# 100
16) Phenanthrene	17.03	178	1761	0.10	ng	97
17) Anthracene	17.11	178	1458	0.10	ng	98
18) Fluoranthene	19.05	202	1758	0.10	ng	99
20) Pyrene	19.42	202	1802	0.10	ng	99
22) Benzo(a)anthracene	21.17	228	1695	0.11	ng	94
23) Chrysene	21.22	228	1637	0.10	ng	95
24) Indeno(1,2,3-cd)pyrene	25.53	276	1671	0.10	ng	100
26) Benzo(b)fluoranthene	22.70	252	1661	0.10	ng	# 76
27) Benzo(k)fluoranthene	22.75	252	1424	0.09	ng	# 73
28) Benzo(a)pyrene	23.25	252	1325	0.10	ng	# 69
29) Dibenzo(a,h)anthracene	25.54	278	1303	0.10	ng	# 78
30) Benzo(a,h,i)perylene	26.18	276	1480	0.10	ng	# 85

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

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