

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061915\
 Data File : BM001747.D
 Acq On : 18 Jun 2015 21:23
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

apatel
 6/22/2015 2:45:38 PM

Quant Time: Jun 19 05:02:22 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 04:18:06 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	15703	5.00	ng	0.00
6) Naphthalene-d8	10.46	136	56931	5.00	ng	0.00
12) Acenaphthene-d10	14.32	164	26815	5.00	ng	0.00
18) Phenanthrene-d10	17.06	188	65818	5.00	ng	0.00
25) Chrysene-d12	21.25	240	39814	5.00	ng	0.00
32) Perylene-d12	23.47	264	34745	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	426	0.11	ng	0.00
5) Phenol-d6	6.83	99	540	0.11	ng	0.00
7) Nitrobenzene-d5	8.83	82	517	0.12	ng	0.00
13) 2,4,6-Tribromophenol	15.80	330	86	0.09	ng	0.00
14) 2-Fluorobiphenyl	12.94	172	1190	0.13	ng	0.00
27) Terphenyl-d14	19.70	244	725	0.12	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.49	42	305	0.11	ng	# 95
8) Nitrobenzene	8.87	77	521	0.11	ng	99
9) Naphthalene	10.49	128	1728	0.12	ng	# 91
10) Hexachlorobutadiene	10.78	225	303	0.13	ng	98
11) 2-Methylnaphthalene	12.12	142	1072	0.12	ng	94
15) Acenaphthylene	14.03	152	1434	0.11	ng	99
16) Acenaphthene	14.37	154	967	0.12	ng	98
19) 4-Bromophenyl-phenylether	16.24	248	314	0.05	ng	94
21) Pentachlorophenol	16.72	266	86	0.06	ng	# 80
22) Phenanthrene	17.09	178	2658	0.12	ng	96
23) Anthracene	17.19	178	1035m	0.04	ng	
24) Fluoranthene	19.12	202	1651	0.06	ng	99
26) Pyrene	19.49	202	1692	0.12	ng	99
28) Benzo(a)anthracene	21.24	228	1467	0.12	ng	93
29) Chrysene	21.28	228	1580	0.13	ng	95
31) Indeno(1,2,3-cd)pyrene	25.70	276	1414	0.14	ng	99
33) Benzo(b)fluoranthene	22.80	252	1391	0.12	ng	# 81
34) Benzo(k)fluoranthene	22.84	252	1232	0.11	ng	# 83
35) Benzo(a)pyrene	23.38	252	1134	0.12	ng	# 74
36) Dibenzo(a,h)anthracene	25.72	278	1131	0.14	ng	# 81
37) Benzo(g,h,i)perylene	26.39	276	1244	0.14	ng	# 85

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

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