

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050115\
 Data File : BM001146.D
 Acq On : 02 May 2015 00:20
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

apatel
 5/5/2015 8:03:34 AM

Quant Time: May 04 16:32:56 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM050115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 02 01:33:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	16258	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	62506	5.00	ng	0.00
12) Acenaphthene-d10	14.20	164	34611	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	83956	5.00	ng	0.00
24) Chrysene-d12	21.15	240	73268	5.00	ng	0.00
30) Perylene-d12	23.29	264	64730	5.00	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.13	112	296	0.09	ng	0.00
5) Phenol-d6	6.72	99	343	0.07	ng	0.02
7) Nitrobenzene-d5	8.70	82	280	0.08	ng	0.01
13) 2,4,6-Tribromophenol	15.71	330	95m	0.06	ng	0.02
14) 2-Fluorobiphenyl	12.81	172	1123	0.10	ng	0.00
26) Terphenyl-d14	19.59	244	938	0.09	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.37	42	175	0.09	ng	# 8
8) Nitrobenzene	8.74	77	289	0.06	ng	96
9) Naphthalene	10.36	128	1350	0.09	ng	# 87
10) Hexachlorobutadiene	10.65	225	269	0.10	ng	99
11) 2-Methylnaphthalene	11.99	142	890	0.09	ng	# 87
15) Acenaphthylene	13.91	152	1273	0.08	ng	# 100
16) Acenaphthene	14.24	154	1116	0.10	ng	90
17) Fluorene	15.24	166	1108	0.09	ng	99
19) Hexachlorobenzene	16.26	284	483	0.10	ng	# 89
21) Phenanthrene	16.98	178	2170	0.10	ng	97
22) Anthracene	17.08	178	1629	0.08	ng	96
23) Fluoranthene	19.01	202	2035	0.09	ng	100
25) Pvrene	19.38	202	2091	0.09	ng	99
27) Benzo(a)anthracene	21.13	228	2042	0.10	ng	97
28) Chrysene	21.18	228	1998	0.09	ng	97
29) Indeno(1,2,3-cd)pvrene	25.43	276	1736	0.08	ng	99
31) Benzo(b)fluoranthene	22.65	252	1699	0.08	ng	# 83
32) Benzo(k)fluoranthene	22.69	252	1844	0.09	ng	# 83
33) Benzo(a)pvrene	23.20	252	1499	0.08	ng	# 76
34) Dibenzo(a,h)anthracene	25.45	278	1346	0.08	ng	# 76
35) Benzo(q,h,i)perylene	26.09	276	1476	0.08	ng	# 86

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

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