

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060615\
 Data File : BM001573.D
 Acq On : 05 Jun 2015 19:37
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

apatel
 6/8/2015 2:47:34 PM

Quant Time: Jun 06 06:15:33 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 06 01:16:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	26638	5.00	ng	0.00
6) Naphthalene-d8	10.47	136	94572	5.00	ng	0.00
12) Acenaphthene-d10	14.35	164	44103	5.00	ng	0.00
18) Phenanthrene-d10	17.09	188	79476	5.00	ng	0.00
25) Chrysene-d12	21.28	240	70918	5.00	ng	0.00
32) Perylene-d12	23.52	264	63024	5.00	ng	0.01

System Monitoring Compounds

4) 2-Fluorophenol	5.28	112	2894	0.48	ng	0.00
5) Phenol-d6	6.86	99	3490	0.47	ng	0.00
7) Nitrobenzene-d5	8.86	82	3037	0.46	ng	0.00
13) 2,4,6-Tribromophenol	15.83	330	858	0.47	ng	0.00
14) 2-Fluorobiphenyl	12.97	172	6801	0.52	ng	0.00
27) Terphenyl-d14	19.72	244	4309	0.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.21	88	1379	0.49	ng	95
3) n-Nitrosodimethylamine	3.52	42	2004	0.49	ng	# 98
8) Nitrobenzene	8.89	77	3233	0.46	ng	99
9) Naphthalene	10.52	128	10055	0.51	ng	99
10) Hexachlorobutadiene	10.82	225	1750	0.51	ng	99
11) 2-Methylnaphthalene	12.15	142	6183	0.49	ng	99
15) Acenaphthylene	14.06	152	9072	0.47	ng	100
16) Acenaphthene	14.41	154	5918	0.51	ng	99
17) Fluorene	15.39	166	3721	0.41	ng	# 100
19) 4-Bromophenyl-phenylether	16.27	248	2448	0.73	ng	98
20) Hexachlorobenzene	16.38	284	1166	0.67	ng	# 72
21) Pentachlorophenol	16.75	266	469	0.51	ng	94
22) Phenanthrene	17.12	178	12866	0.55	ng	99
23) Anthracene	17.23	178	5782m	0.54	ng	
24) Fluoranthene	19.15	202	10116	0.60	ng	99
26) Pyrene	19.52	202	10751	0.49	ng	100
28) Benzo(a)anthracene	21.27	228	9377	0.47	ng	# 85
29) Chrysene	21.31	228	9425	0.53	ng	97
30) Bis(2-ethylhexyl)phthalate	21.19	149	4795	0.34	ng	100
31) Indeno(1,2,3-cd)pyrene	25.77	276	9813	0.50	ng	99
33) Benzo(b)fluoranthene	22.84	252	9219	0.53	ng	97
34) Benzo(k)fluoranthene	22.89	252	8213	0.50	ng	98
35) Benzo(a)pyrene	23.42	252	7893	0.50	ng	97
36) Dibenzo(a,h)anthracene	25.79	278	7721	0.52	ng	97
37) Benzo(g,h,i)perylene	26.46	276	8227	0.53	ng	98

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

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