

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060315\
 Data File : BM001540.D
 Acq On : 03 Jun 2015 18:24
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

apatel
 6/5/2015 8:04:00 AM

Quant Time: Jun 04 03:59:46 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 03:14:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	29178	5.00	ng	0.00
6) Naphthalene-d8	10.49	136	104831	5.00	ng	0.00
12) Acenaphthene-d10	14.35	164	49962	5.00	ng	0.00
18) Phenanthrene-d10	17.09	188	126260	5.00	ng	0.00
25) Chrysene-d12	21.28	240	79060	5.00	ng	0.00
32) Perylene-d12	23.53	264	77779	5.00	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.28	112	65256	10.16	ng	0.00
5) Phenol-d6	6.86	99	82279	10.37	ng	0.00
7) Nitrobenzene-d5	8.86	82	76328	11.49	ng	0.00
13) 2,4,6-Tribromophenol	15.83	330	20616	19.26	ng	0.00
14) 2-Fluorobiphenyl	12.97	172	142296	9.51	ng	0.00
27) Terphenyl-d14	19.73	244	97081	9.93	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.21	88	25462	8.35	ng	96
3) n-Nitrosodimethylamine	3.51	42	43512	10.46	ng	97
8) Nitrobenzene	8.90	77	80892	11.47	ng	99
9) Naphthalene	10.52	128	210067	9.50	ng	98
10) Hexachlorobutadiene	10.82	225	36008	9.34	ng	100
11) 2-Methylnaphthalene	12.16	142	134088	9.65	ng	99
15) Acenaphthylene	14.07	152	225491	10.53	ng	100
16) Acenaphthene	14.41	154	131394	9.90	ng	95
17) Fluorene	15.39	166	118175	5.67	ng	99
19) 4-Bromophenyl-phenylether	16.27	248	40126	18.55	ng	# 88
20) Hexachlorobenzene	16.41	284	24132	5.39	ng	# 100
21) Pentachlorophenol	16.75	266	16616	8.49	ng	95
22) Phenanthrene	17.12	178	327380	13.62	ng	99
23) Anthracene	17.23	178	157525m	6.19	ng	
24) Fluoranthene	19.15	202	234892	10.35	ng	99
26) Pyrene	19.52	202	244408	10.18	ng	100
28) Benzo(a)anthracene	21.27	228	219826	10.27	ng	97
29) Chrysene	21.31	228	190835	9.46	ng	97
30) Bis(2-ethylhexyl)phthalate	21.21	149	178482	11.66	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	221416	9.92	ng	99
33) Benzo(b)fluoranthene	22.85	252	217070	9.86	ng	96
34) Benzo(k)fluoranthene	22.90	252	180515	9.20	ng	97
35) Benzo(a)pyrene	23.43	252	187287	9.85	ng	97
36) Dibenzo(a,h)anthracene	25.80	278	173140	9.23	ng	95
37) Benzo(g,h,i)perylene	26.48	276	179359	9.14	ng	99

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

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