

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
 Data File : BM001651.D
 Acq On : 11 Jun 2015 20:58
 Operator : TP/IZ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

apatel
 6/15/2015 12:41:12 PM

Quant Time: Jun 12 16:35:34 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 02:07:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	20711	5.00	ng	0.00
6) Naphthalene-d8	10.47	136	78720	5.00	ng	0.00
12) Acenaphthene-d10	14.34	164	39568	5.00	ng	0.00
18) Phenanthrene-d10	17.06	188	74206	5.00	ng	0.00
25) Chrysene-d12	21.28	240	70340	5.00	ng	0.01
32) Perylene-d12	23.51	264	55322	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.27	112	23537	4.99	ng	0.00
5) Phenol-d6	6.84	99	30733	5.28	ng	0.00
7) Nitrobenzene-d5	8.85	82	30059	5.38	ng	0.00
13) 2,4,6-Tribromophenol	15.84	330	6049	3.82	ng	0.00
14) 2-Fluorobiphenyl	12.95	172	62815	4.86	ng	0.00
27) Terphenyl-d14	19.72	244	51555	5.50	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.21	88	10440	4.55	ng	97
3) n-Nitrosodimethylamine	3.51	42	17573	5.32	ng	# 98
8) Nitrobenzene	8.89	77	33244	5.53	ng	98
9) Naphthalene	10.52	128	86225	4.86	ng	99
10) Hexachlorobutadiene	10.82	225	14277	4.72	ng	99
11) 2-Methylnaphthalene	12.14	142	57296	5.23	ng	94
15) Acenaphthylene	14.05	152	94373	5.30	ng	100
16) Acenaphthene	14.40	154	54035	4.84	ng	98
20) Hexachlorobenzene	16.38	284	29982	6.91	ng	# 55
21) Pentachlorophenol	16.72	266	11624	8.61	ng	90
23) Anthracene	17.20	178	142107m	7.93	ng	
26) Pyrene	19.51	202	117995	5.13	ng	100
28) Benzo(a)anthracene	21.25	228	104027	5.01	ng	95
29) Chrysene	21.31	228	97357m	4.97	ng	
30) Bis(2-ethylhexyl)phthalate	21.19	149	74929	6.21	ng	98
31) Indeno(1,2,3-cd)pyrene	25.76	276	76111	3.97	ng	100
33) Benzo(b)fluoranthene	22.83	252	88287	5.27	ng	98
34) Benzo(k)fluoranthene	22.87	252	83109	5.14	ng	98
35) Benzo(a)pyrene	23.40	252	74323	5.06	ng	98
36) Dibenzo(a,h)anthracene	25.77	278	59864	4.52	ng	98
37) Benzo(g,h,i)perylene	26.45	276	61628	4.32	ng	99

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

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