

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
 Data File : BM001677.D
 Acq On : 12 Jun 2015 19:42
 Operator : TP/IZ
 Sample : SSTDIC010
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDIC010

Quant Time: Jun 13 00:27:10 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 13 00:26:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	14161	5.00	ng	0.00
6) Naphthalene-d8	10.46	136	54543	5.00	ng	0.00
12) Acenaphthene-d10	14.32	164	27709	5.00	ng	0.00
18) Phenanthrene-d10	17.06	188	87408	5.00	ng	0.00
25) Chrysene-d12	21.25	240	52161	5.00	ng	0.00
32) Perylene-d12	23.48	264	47992	5.00	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.25	112	34163	10.13	ng	0.00
5) Phenol-d6	6.83	99	45483	10.71	ng	0.00
7) Nitrobenzene-d5	8.83	82	45008	10.68	ng	0.00
13) 2,4,6-Tribromophenol	15.80	330	6929	7.29	ng	0.00
14) 2-Fluorobiphenyl	12.94	172	92112	9.44	ng	0.00
27) Terphenyl-d14	19.70	244	74978	9.51	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.18	88	14951	8.41	ng	98
3) n-Nitrosodimethylamine	3.48	42	26520	10.69	ng	94
8) Nitrobenzene	8.87	77	49230	10.80	ng	98
9) Naphthalene	10.51	128	123058	9.25	ng	98
10) Hexachlorobutadiene	10.78	225	20562	9.27	ng	99
11) 2-Methylnaphthalene	12.13	142	83179	9.71	ng	100
15) Acenaphthylene	14.03	152	140850	10.80	ng	100
16) Acenaphthene	14.39	154	78958	9.52	ng	99
17) Fluorene	15.36	166	125859	13.98	ng	99
19) 4-Bromophenyl-phenylether	16.24	248	15768	2.08	ng	89
20) Hexachlorobenzene	16.38	284	29803	4.07	ng	# 100
21) Pentachlorophenol	16.72	266	20802	9.94	ng	98
22) Phenanthrene	17.09	178	194935	7.03	ng	98
23) Anthracene	17.19	178	166261	5.28	ng	99
24) Fluoranthene	19.13	202	164483	4.31	ng	100
26) Pyrene	19.49	202	174565	9.74	ng	99
28) Benzo(a)anthracene	21.24	228	156724	9.54	ng	97
29) Chrysene	21.28	228	146768	9.33	ng	97
30) Bis(2-ethylhexyl)phthalate	21.18	149	117302	11.56	ng	99
31) Indeno(1,2,3-cd)pyrene	25.72	276	163739	12.67	ng	100
33) Benzo(b)fluoranthene	22.81	252	152546	9.76	ng	97
34) Benzo(k)fluoranthene	22.85	252	138343	9.09	ng	96
35) Benzo(a)pyrene	23.39	252	135287	10.30	ng	97
36) Dibenzo(a,h)anthracene	25.74	278	132266	11.67	ng	96
37) Benzo(g,h,i)perylene	26.41	276	133759	10.97	ng	98

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

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