Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE022615\
Data File : BE089706.D
Acq On : 27 Feb 2015 1:22
Operator : TP/JJ
Sample : SSTDICC005
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC005

Quant Time: Feb 27 04:58:02 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE022615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 27 04:47:17 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.33	152	54773	5.00	ng		0.00
7) Naphthalene-d8	12.23	136	229646	5.00	ng		0.00
13) Acenaphthene-d10	16.41	164	105922	5.00	ng		0.00
17) Phenanthrene-d10	19.76	188	194465	5.00	ng		0.00
20) Chrysene-d12	23.65	240	176532	5.00	ng		0.00
27) Perylene-d12	25.33	264	138313	5.00	ng		0.00
System Monitoring Compounds							
3) 2-Fluorophenol	6.49	112	72319	4.32	ng		0.00
4) Phenol-d6	8.61	99	95640	4.61	ng		0.00
8) Nitrobenzene-d5	10.60	82	68192	4.36	ng		0.00
14) 2,4,6-Tribromophenol	18.31	330	7867	3.57	ng		0.00
15) 2-Fluorobiphenyl	14.87	172	116349	3.73	ng		0.00
22) Terphenyl-d14	22.30	244	110913	5.70	ng		0.00
Target Compounds							Qvalue
2) n-Nitrosodimethylamine	3.32	42	45636	5.14	ng	#	81
5) Phenol	8.63	94	103833	4.70	ng	#	25
6) bis(2-Chloroethyl)ether	8.82	93	78413	4.75	ng	#	100
9) Nitrobenzene	10.64	77	81318	4.59	ng		98
10) 2,4-Dimethylphenol	11.57	122	60002	4.43	ng		98
11) 2,4-Dichlorophenol	11.95	162	45929	4.44	ng		98
12) Hexachlorobutadiene	12.62	225	22337	3.08	ng		100
16) 2,4-Dinitrophenol	16.65	184	2762	2.82	ng	#	75
18) Hexachlorobenzene	18.99	284	25872	3.04	ng	#	83
19) Pentachlorophenol	19.43	266	7140	3.22	ng		92
21) Benzidine	21.95	184	68739	5.02	ng		97
23) Benzo(a)anthracene	23.64	228	714972	4.52	ng		93
24) 3,3'-Dichlorobenzidine	23.63	252	49723	2.88	ng	#	96
25) Chrysene	23.68	228	761057	3.92	ng		97
26) Indeno(1,2,3-cd)pyrene	26.74	276	135805	2.60	ng	#	76
28) Benzo(b)fluoranthene	24.91	252	120314	5.01	ng		98
29) Benzo(k)fluoranthene	24.94	252	187618	4.86	ng		99
30) Benzo(a)pyrene	25.27	252	118763	3.79	ng		99
31) Dibenzo(a,h)anthracene	26.77	278	105719	3.17	ng		99

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

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