Quantitation Report (Not Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
Data File : BE089425.D
Acq On : 29 Jan 2015 23:01
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC001

Manual Integrations APPROVED
apatel
1/30/2015 5:56:01 PM

Quant Time: Jan 30 17:16:30 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 30 17:07:09 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	156924	5.00	ng		0.00
7) Naphthalene-d8	12.50	136	595883	5.00	ng		0.00
13) Acenaphthene-d10	16.68	164	244350	5.00	ng		0.00
17) Phenanthrene-d10	19.98	188	491333	5.00	ng		0.00
20) Chrysene-d12	23.84	240	601211	5.00	ng		0.00
27) Perylene-d12	25.55	264	632987	5.00	ng		0.00
System Monitoring Compounds							
3) 2-Fluorophenol	6.74	112	53927	0.91	ng		0.00
4) Phenol-d6	8.85	99	65603	0.94	ng		0.00
8) Nitrobenzene-d5	10.84	82	46062	0.95	ng		0.00
14) 2,4,6-Tribromophenol	18.58	330	6759	1.00	ng		0.00
15) 2-Fluorobiphenyl	15.13	172	64133	0.94	ng		0.00
22) Terphenyl-d14	22.48	244	76027	1.09	ng		0.00
Target Compounds							Qvalue
2) n-Nitrosodimethylamine	3.58	42	30442	0.92	ng	#	58
5) Phenol	8.88	94	72194	0.94	ng		86
6) bis(2-Chloroethyl)ether	9.06	93	53404	0.93	ng	#	99
9) Nitrobenzene	10.88	77	52983	0.96	ng		99
10) 2,4-Dimethylphenol	11.82	122	37292	0.95	ng		93
11) 2,4-Dichlorophenol	12.21	162	26571	0.97	ng		93
12) Hexachlorobutadiene	12.89	225	17131	0.93	ng		99
16) 2,4-Dinitrophenol	16.89	184	1813	1.26	ng		93
18) Hexachlorobenzene	19.23	284	23191	0.96	ng		96
19) Pentachlorophenol	19.65	266	3661	0.86	ng		97
21) Benzidine	22.14	184	58307	1.27	ng		99
23) Benzo(a)anthracene	23.83	228	479971	1.05	ng		100
24) 3,3'-Dichlorobenzidine	23.81	252	39679	1.03	ng		97
25) Chrysene	23.87	228	579205	0.93	ng		97
26) Indeno(1,2,3-cd)pyrene	27.07	276	121984	1.09	ng	#	100
28) Benzo(b)fluoranthene	25.13	252	258496	3.39	ng		99
29) Benzo(k)fluoranthene	25.13	252	260667	1.27	ng		100
30) Benzo(a)pyrene	25.48	252	85235	1.09	ng		99
31) Dibenzo(a,h)anthracene	27.09	278	93960	1.04	ng		97

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

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