Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
Data File : BE089413.D
Acq On : 29 Jan 2015 13:19
Operator : TP/IZ
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDICCC001

Manual Integrations APPROVED
apatel
1/30/2015 5:55:33 PM

Quant Time: Jan 30 02:59:58 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 02:34:28 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	172715	5.00	ng		0.00
7) Naphthalene-d8	12.50	136	627822	5.00	ng		0.00
13) Acenaphthene-d10	16.68	164	239675	5.00	ng		0.00
17) Phenanthrene-d10	19.99	188	479216	5.00	ng		0.00
20) Chrysene-d12	23.84	240	648652	5.00	ng		0.00
27) Perylene-d12	25.55	264	691358	5.00	ng		0.00
System Monitoring Compounds							
3) 2-Fluorophenol	6.74	112	61007	1.10	ng		0.00
4) Phenol-d6	8.85	99	72163	1.06	ng		0.00
8) Nitrobenzene-d5	10.84	82	48746	1.08	ng		0.00
14) 2,4,6-Tribromophenol	18.57	330	6328	1.14	ng		0.00
15) 2-Fluorobiphenyl	15.13	172	64150	0.89	ng		0.00
22) Terphenyl-d14	22.48	244	74711	1.04	ng		0.00
Target Compounds							Qvalue
2) n-Nitrosodimethylamine	3.58	42	35276	1.20	ng		100
5) Phenol	8.87	94	78953	1.09	ng		100
6) bis(2-Chloroethyl)ether	9.06	93	58655	1.09	ng	#	100
9) Nitrobenzene	10.89	77	55865	1.09	ng		100
10) 2,4-Dimethylphenol	11.82	122	39111	1.02	ng		100
11) 2,4-Dichlorophenol	12.21	162	27808	0.95	ng		100
12) Hexachlorobutadiene	12.89	225	18147	0.88	ng		100
16) 2,4-Dinitrophenol	16.89	184	1173	0.72	ng		100
18) Hexachlorobenzene	19.23	284	22085	0.99	ng		100
19) Pentachlorophenol	19.65	266	3679	0.86	ng		100
21) Benzidine	22.14	184	37078	0.71	ng		100
23) Benzo(a)anthracene	23.83	228	472429	0.80	ng		100
24) 3,3'-Dichlorobenzidine	23.81	252	35384	0.75	ng		100
25) Chrysene	23.87	228	631594	0.87	ng		100
26) Indeno(1,2,3-cd)pyrene	27.06	276	103305	0.62	ng	#	100
28) Benzo(b)fluoranthene	25.10	252	56283m	0.64	ng		
29) Benzo(k)fluoranthene	25.13	252	200329m	1.01	ng		
30) Benzo(a)pyrene	25.48	252	75116	0.71	ng		100
31) Dibenzo(a,h)anthracene	27.08	278	74833	0.69	ng		100

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

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