Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
Data File : BE089471.D
Acq On : 3 Feb 2015 19:36
Operator : TP/IZ
Sample : SSTDICC0.2
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDICC0.2

Manual Integrations APPROVED
apatel
2/4/2015 2:55:09 PM

Quant Time: Feb 04 00:44:11 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE020315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 04 00:22:17 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.53	152	57169	5.00	ng		0.00
7) Naphthalene-d8	12.43	136	225647	5.00	ng		0.00
13) Acenaphthene-d10	16.61	164	96710	5.00	ng		0.00
17) Phenanthrene-d10	19.93	188	186273	5.00	ng		0.00
20) Chrysene-d12	23.78	240	180368	5.00	ng		0.00
27) Perylene-d12	25.49	264	170057	5.00	ng		0.00
System Monitoring Compounds							
3) 2-Fluorophenol	6.72	112	3829	0.20	ng		0.00
4) Phenol-d6	8.83	99	4834	0.22	ng		0.00
8) Nitrobenzene-d5	10.79	82	3377	0.21	ng		0.00
14) 2,4,6-Tribromophenol	18.51	330	569	0.23	ng		0.00
15) 2-Fluorobiphenyl	15.06	172	5245	0.19	ng		0.00
22) Terphenyl-d14	22.43	244	5031	0.24	ng		0.00
Target Compounds							Qvalue
2) n-Nitrosodimethylamine	3.60	42	1863	0.20	ng	#	76
5) Phenol	8.86	94	5171m	0.21	ng		
6) bis(2-Chloroethyl)ether	9.01	93	3946	0.22	ng	#	99
9) Nitrobenzene	10.83	77	3834	0.21	ng		99
10) 2,4-Dimethylphenol	11.77	122	2797	0.21	ng		97
11) 2,4-Dichlorophenol	12.15	162	2034	0.20	ng		94
12) Hexachlorobutadiene	12.81	225	1530	0.22	ng		96
18) Hexachlorobenzene	19.16	284	1943	0.20	ng	#	71
19) Pentachlorophenol	19.60	266	97	0.08	ng	#	64
21) Benzidine	22.09	184	927	0.07	ng	#	83
23) Benzo(a)anthracene	23.77	228	31508	0.22	ng		98
24) 3,3'-Dichlorobenzidine	23.76	252	2092	0.16	ng	#	95
25) Chrysene	23.82	228	35190	0.20	ng		97
26) Indeno(1,2,3-cd)pyrene	26.96	276	5374	0.11	ng	#	86
28) Benzo(b)fluoranthene	25.04	252	3773m	0.16	ng		
29) Benzo(k)fluoranthene	25.07	252	8639m	0.20	ng		
30) Benzo(a)pyrene	25.42	252	4123	0.16	ng	#	90
31) Dibenzo(a,h)anthracene	26.99	278	4464m	0.14	ng		

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

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