

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089468.D
 Acq On : 3 Feb 2015 17:37
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICCC001

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:54:57 PM

Quant Time: Feb 04 00:38:42 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	59878	5.00	ng	0.00
7) Naphthalene-d8	12.43	136	239834	5.00	ng	0.00
13) Acenaphthene-d10	16.61	164	102742	5.00	ng	0.00
17) Phenanthrene-d10	19.93	188	197851	5.00	ng	0.00
20) Chrysene-d12	23.78	240	198084	5.00	ng	0.00
27) Perylene-d12	25.49	264	188603	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.72	112	18812	0.94	ng	0.00
4) Phenol-d6	8.84	99	23935	1.02	ng	0.00
8) Nitrobenzene-d5	10.79	82	17607	1.01	ng	0.00
14) 2,4,6-Tribromophenol	18.51	330	3456	1.29	ng	0.00
15) 2-Fluorobiphenyl	15.06	172	26757	0.93	ng	0.00
22) Terphenyl-d14	22.43	244	28053	1.23	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.55	42	10838	1.13	ng	100
5) Phenol	8.86	94	26282m	1.04	ng	
6) bis(2-Chloroethyl)ether	9.01	93	18794	1.02	ng	# 100
9) Nitrobenzene	10.83	77	20424	1.03	ng	100
10) 2,4-Dimethylphenol	11.77	122	14751	1.04	ng	100
11) 2,4-Dichlorophenol	12.15	162	10871	1.01	ng	100
12) Hexachlorobutadiene	12.81	225	7304	0.97	ng	100
16) 2,4-Dinitrophenol	16.84	184	344	0.56	ng	100
18) Hexachlorobenzene	19.17	284	9996	0.96	ng	100
19) Pentachlorophenol	19.61	266	1224	1.00	ng	100
21) Benzidine	22.09	184	9469	0.67	ng	100
23) Benzo(a)anthracene	23.77	228	178737	1.12	ng	100
24) 3,3'-Dichlorobenzidine	23.76	252	14179	1.01	ng	100
25) Chrysene	23.82	228	178468	0.93	ng	100
26) Indeno(1,2,3-cd)pyrene	26.97	276	43391	0.77	ng	100
28) Benzo(b)fluoranthene	25.04	252	32536	1.26	ng	100
29) Benzo(k)fluoranthene	25.07	252	52741	1.11	ng	100
30) Benzo(a)pyrene	25.42	252	33659	1.17	ng	100
31) Dibenzo(a,h)anthracene	26.99	278	37166	1.07	ng	100

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

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