

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021615\
 Data File : BE089574.D
 Acq On : 16 Feb 2015 17:50
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

apatel
 2/17/2015 3:36:17 PM

Quant Time: Feb 16 18:30:50 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIM-BE021615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 16 17:44:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	77828	5.00	ng	0.00
7) Naphthalene-d8	12.25	136	307762	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	144300	5.00	ng	0.00
17) Phenanthrene-d10	19.77	188	276447	5.00	ng	0.00
20) Chrysene-d12	23.66	240	306517	5.00	ng	0.00
27) Perylene-d12	25.34	264	292141	5.00	ng	0.00

System Monitoring Compounds

3) 2-Fluorophenol	6.50	112	10445	0.46	ng	0.00
4) Phenol-d6	8.62	99	13911	0.46	ng	0.00
8) Nitrobenzene-d5	10.61	82	5645	0.42	ng	0.00
14) 2,4,6-Tribromophenol	18.31	330	1371	0.41	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	19119	0.49	ng	0.00
22) Terphenyl-d14	22.30	244	19591	0.46	ng	0.00

Target Compounds

						Qvalue
2) n-Nitrosodimethylamine	3.39	42	5051m	0.39	ng	
5) Phenol	8.65	94	14984	0.46	ng	93
6) bis(2-Chloroethyl)ether	8.83	93	11585	0.47	ng	# 98
9) Nitrobenzene	10.65	77	6353	0.39	ng	100
10) 2,4-Dimethylphenol	11.59	122	8130	0.44	ng	100
11) 2,4-Dichlorophenol	11.97	162	6487	0.45	ng	99
12) Hexachlorobutadiene	12.64	225	4643	0.50	ng	99
16) 2,4-Dinitrophenol	16.67	184	128	0.23	ng	# 68
18) Hexachlorobenzene	19.00	284	5791	0.50	ng	99
19) Pentachlorophenol	19.45	266	768	0.35	ng	94
21) Benzidine	21.96	184	7923	0.27	ng	98
23) Benzo(a)anthracene	23.64	228	135904	0.44	ng	97
24) 3,3'-Dichlorobenzidine	23.64	252	7539	0.32	ng	99
25) Chrysene	23.69	228	143644	0.49	ng	100
26) Indeno(1,2,3-cd)pyrene	26.75	276	30816	0.39	ng	# 81
28) Benzo(b)fluoranthene	24.91	252	23187	0.35	ng	100
29) Benzo(k)fluoranthene	24.94	252	38401	0.46	ng	99
30) Benzo(a)pyrene	25.28	252	23008	0.36	ng	99
31) Dibenzo(a,h)anthracene	26.78	278	23759	0.37	ng	100

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

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