

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021615\
 Data File : BE089572.D
 Acq On : 16 Feb 2015 16:27
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICCC001

Manual Integrations
 APPROVED

apatel
 2/17/2015 3:36:09 PM

Quant Time: Feb 16 17:25:36 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 17:13:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	78577	5.00	ng	0.00
7) Naphthalene-d8	12.25	136	312599	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	144642	5.00	ng	0.00
17) Phenanthrene-d10	19.77	188	273646	5.00	ng	0.00
20) Chrysene-d12	23.65	240	310751	5.00	ng	0.00
27) Perylene-d12	25.34	264	294199	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	6.50	112	22740	0.88	ng	0.00
4) Phenol-d6	8.63	99	29913	0.93	ng	0.00
8) Nitrobenzene-d5	10.61	82	12705	0.58	ng	0.00
14) 2,4,6-Tribromophenol	18.31	330	3080	1.05	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	40391	0.90	ng	0.00
22) Terphenyl-d14	22.30	244	42294	1.36	ng	0.00

Target Compounds				Qvalue		
2) n-Nitrosodimethylamine	3.36	42	12082m	0.92	ng	
5) Phenol	8.65	94	32522	0.99	ng	100
6) bis(2-Chloroethyl)ether	8.83	93	24640	1.00	ng	# 100
9) Nitrobenzene	10.65	77	14968	0.60	ng	100
10) 2,4-Dimethylphenol	11.59	122	18435	0.85	ng	100
11) 2,4-Dichlorophenol	11.97	162	14431	0.93	ng	100
12) Hexachlorobutadiene	12.64	225	9672	0.95	ng	100
16) 2,4-Dinitrophenol	16.66	184	374	0.42	ng	100
18) Hexachlorobenzene	19.00	284	11770	0.88	ng	100
19) Pentachlorophenol	19.44	266	1818	0.74	ng	100
21) Benzidine	21.96	184	25255	0.97	ng	100
23) Benzo(a)anthracene	23.64	228	303353	1.30	ng	100
24) 3,3'-Dichlorobenzidine	23.64	252	20597	0.97	ng	100
25) Chrysene	23.69	228	297861	0.83	ng	100
26) Indeno(1,2,3-cd)pyrene	26.74	276	74074	0.45	ng	100
28) Benzo(b)fluoranthene	24.91	252	58499	1.88	ng	100
29) Benzo(k)fluoranthene	24.94	252	82217	1.31	ng	100
30) Benzo(a)pyrene	25.28	252	57199	1.55	ng	100
31) Dibenzo(a,h)anthracene	26.77	278	58609	1.73	ng	100

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

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