

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE022615\
 Data File : BE089704.D
 Acq On : 27 Feb 2015 00:04
 Operator : TP/JJ
 Sample : SSTDICC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICCC001

Manual Integrations
 APPROVED

apatel
 2/27/2015 4:13:06 PM

Quant Time: Feb 27 05:06:43 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE022615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 27 04:47:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.33	152	62553	5.00	ng	0.00
7) Naphthalene-d8	12.23	136	272690	5.00	ng	0.00
13) Acenaphthene-d10	16.41	164	126771	5.00	ng	0.00
17) Phenanthrene-d10	19.76	188	228734	5.00	ng	0.00
20) Chrysene-d12	23.65	240	184375	5.00	ng	0.00
27) Perylene-d12	25.33	264	144099	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.49	112	14149	0.74	ng	0.00
4) Phenol-d6	8.61	99	19189	0.81	ng	0.00
8) Nitrobenzene-d5	10.60	82	12758	0.69	ng	0.00
14) 2,4,6-Tribromophenol	18.30	330	1009	0.38	ng	0.00
15) 2-Fluorobiphenyl	14.87	172	26280	0.70	ng	0.00
22) Terphenyl-d14	22.30	244	20528	1.01	ng	0.00
Target Compounds						
2) n-Nitrosodimethylamine	3.36	42	8876m	0.88	ng	Qvalue
5) Phenol	8.64	94	21031	0.83	ng	# 57
6) bis(2-Chloroethyl)ether	8.82	93	16362	0.87	ng	# 100
9) Nitrobenzene	10.64	77	15225	0.72	ng	100
10) 2,4-Dimethylphenol	11.57	122	11612	0.72	ng	100
11) 2,4-Dichlorophenol	11.95	162	8422	0.69	ng	100
12) Hexachlorobutadiene	12.62	225	4748	0.55	ng	100
16) 2,4-Dinitrophenol	16.65	184	294	0.84	ng	100
18) Hexachlorobenzene	18.98	284	5452	0.55	ng	100
19) Pentachlorophenol	19.43	266	668	0.55	ng	100
21) Benzidine	21.96	184	6086	0.43	ng	100
23) Benzo(a)anthracene	23.64	228	117331	0.71	ng	100
24) 3,3'-Dichlorobenzidine	23.63	252	5264	0.61	ng	100
25) Chrysene	23.68	228	146318	0.72	ng	100
26) Indeno(1,2,3-cd)pyrene	26.74	276	16125	0.54	ng	# 48
28) Benzo(b)fluoranthene	24.91	252	11410	0.67	ng	100
29) Benzo(k)fluoranthene	24.94	252	33915	0.84	ng	100
30) Benzo(a)pyrene	25.27	252	13628	0.73	ng	100
31) Dibenzo(a,h)anthracene	26.77	278	11624	0.68	ng	100

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

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