

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089480.D
 Acq On : 4 Feb 2015 1:25
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:55:30 PM

Quant Time: Feb 04 03:04:16 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 00:53:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.53	152	57597	5.00	ng	0.00
7) Naphthalene-d8	12.43	136	230796	5.00	ng	0.00
13) Acenaphthene-d10	16.61	164	101668	5.00	ng	0.00
17) Phenanthrene-d10	19.93	188	191950	5.00	ng	0.00
20) Chrysene-d12	23.78	240	184352	5.00	ng	0.00
27) Perylene-d12	25.49	264	170287	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.73	112	17870	0.96	ng	0.00
4) Phenol-d6	8.85	99	22681	0.97	ng	0.00
8) Nitrobenzene-d5	10.79	82	17013	0.97	ng	0.00
14) 2,4,6-Tribromophenol	18.52	330	3181	0.89	ng	0.00
15) 2-Fluorobiphenyl	15.06	172	26751	1.00	ng	0.00
22) Terphenyl-d14	22.43	244	27560	1.05	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.56	42	10207	0.95	ng	# 95
5) Phenol	8.87	94	24776m	0.97	ng	
6) bis(2-Chloroethyl)ether	9.01	93	17902	0.97	ng	# 100
9) Nitrobenzene	10.83	77	19733	0.99	ng	100
10) 2,4-Dimethylphenol	11.78	122	14161	0.98	ng	100
11) 2,4-Dichlorophenol	12.16	162	10424	0.98	ng	95
12) Hexachlorobutadiene	12.81	225	7133	1.00	ng	99
16) 2,4-Dinitrophenol	16.84	184	238	0.77	ng	93
18) Hexachlorobenzene	19.17	284	10028	1.01	ng	95
19) Pentachlorophenol	19.61	266	1065	0.86	ng	94
21) Benzidine	22.10	184	15144	1.72	ng	98
23) Benzo(a)anthracene	23.77	228	168894	1.03	ng	100
24) 3,3'-Dichlorobenzidine	23.77	252	14472	1.04	ng	# 90
25) Chrysene	23.82	228	168092	0.98	ng	100
26) Indeno(1,2,3-cd)pyrene	26.97	276	44326	1.04	ng	98
28) Benzo(b)fluoranthene	25.04	252	34248	1.04	ng	99
29) Benzo(k)fluoranthene	25.07	252	45362	1.03	ng	99
30) Benzo(a)pyrene	25.42	252	32821	1.01	ng	99
31) Dibenzo(a,h)anthracene	27.00	278	36854	1.05	ng	98

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

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