

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089490.D
 Acq On : 4 Feb 2015 10:07
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:55:47 PM

Quant Time: Feb 04 10:52:41 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	9.54	152	58129	5.00	ng	0.00
7) Naphthalene-d8	12.44	136	231224	5.00	ng	0.00
13) Acenaphthene-d10	16.62	164	99913	5.00	ng	0.00
17) Phenanthrene-d10	19.94	188	186876	5.00	ng	0.00
20) Chrysene-d12	23.79	240	173168	5.00	ng	0.00
27) Perylene-d12	25.49	264	168773	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.76	112	18232	0.97	ng	0.04
4) Phenol-d6	8.89	99	22969	0.97	ng	0.05
8) Nitrobenzene-d5	10.80	82	17655	1.01	ng	0.02
14) 2,4,6-Tribromophenol	18.53	330	3412	0.95	ng	0.02
15) 2-Fluorobiphenyl	15.07	172	27626	1.05	ng	0.00
22) Terphenyl-d14	22.43	244	28378	1.16	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.57	42	10348	0.95	ng	# 75
5) Phenol	8.92	94	22567m	0.87	ng	
6) bis(2-Chloroethyl)ether	9.01	93	17849	0.96	ng	# 99
9) Nitrobenzene	10.84	77	19523	0.98	ng	99
10) 2,4-Dimethylphenol	11.80	122	14172	0.98	ng	96
11) 2,4-Dichlorophenol	12.18	162	10726	1.00	ng	98
12) Hexachlorobutadiene	12.81	225	7187	1.01	ng	98
16) 2,4-Dinitrophenol	16.87	184	221	0.75	ng	94
18) Hexachlorobenzene	19.18	284	9678	1.00	ng	94
19) Pentachlorophenol	19.62	266	727	0.71	ng	94
21) Benzidine	22.11	184	6394	0.82	ng	97
23) Benzo(a)anthracene	23.77	228	161535	1.05	ng	94
24) 3,3'-Dichlorobenzidine	23.77	252	13699	1.05	ng	98
25) Chrysene	23.83	228	157859	0.98	ng	96
26) Indeno(1,2,3-cd)pyrene	26.98	276	46637	1.15	ng	99
28) Benzo(b)fluoranthene	25.05	252	35479	1.08	ng	99
29) Benzo(k)fluoranthene	25.08	252	43635	1.00	ng	99
30) Benzo(a)pyrene	25.43	252	35323	1.09	ng	99
31) Dibenzo(a,h)anthracene	27.01	278	39346	1.12	ng	98

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

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