

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
 Data File : BE089415.D
 Acq On : 29 Jan 2015 14:40
 Operator : TP/IZ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC005

Quant Time: Jan 30 02:45:25 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 30 02:34:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	101241	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	353948	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	124753	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	244053	5.00	ng	0.00
20) Chrysene-d12	23.84	240	398056	5.00	ng	0.00
27) Perylene-d12	25.55	264	451268	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	6.73	112	184444	5.68	ng	0.00
4) Phenol-d6	8.85	99	211290	5.32	ng	0.00
8) Nitrobenzene-d5	10.84	82	142737	5.60	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	18555	6.40	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	171963	4.57	ng	0.00
22) Terphenyl-d14	22.48	244	219573	4.99	ng	0.00

Target Compounds				Qvalue		
2) n-Nitrosodimethylamine	3.55	42	113321	6.57	ng	# 85
5) Phenol	8.87	94	231796	5.48	ng	99
6) bis(2-Chloroethyl)ether	9.06	93	170380	5.41	ng	# 100
9) Nitrobenzene	10.89	77	164018	5.70	ng	99
10) 2,4-Dimethylphenol	11.82	122	112545	5.19	ng	98
11) 2,4-Dichlorophenol	12.21	162	79093	4.79	ng	98
12) Hexachlorobutadiene	12.89	225	52059	4.46	ng	99
16) 2,4-Dinitrophenol	16.89	184	6788	8.03	ng	91
18) Hexachlorobenzene	19.23	284	56340	4.96	ng	94
19) Pentachlorophenol	19.65	266	17030	7.80	ng	98
21) Benzidine	22.14	184	161541	5.01	ng	99
23) Benzo(a)anthracene	23.83	228	1585696	4.39	ng	97
24) 3,3'-Dichlorobenzidine	23.81	252	155264	5.39	ng	98
25) Chrysene	23.87	228	1985069	4.45	ng	99
26) Indeno(1,2,3-cd)pyrene	27.07	276	569746	5.53	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	311357	5.45	ng	99
29) Benzo(k)fluoranthene	25.13	252	693900	5.36	ng	99
30) Benzo(a)pyrene	25.48	252	403000	5.81	ng	99
31) Dibenzo(a,h)anthracene	27.09	278	439366	6.20	ng	98

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

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