

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
 Data File : BE089416.D
 Acq On : 29 Jan 2015 15:19
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC010

Quant Time: Jan 30 02:45:48 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	98013	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	331544	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	113418	5.00	ng	0.00
17) Phenanthrene-d10	19.98	188	221484	5.00	ng	0.00
20) Chrysene-d12	23.84	240	389397	5.00	ng	0.00
27) Perylene-d12	25.55	264	479544	5.00	ng	0.00

System Monitoring Compounds

3) 2-Fluorophenol	6.73	112	354847	11.29	ng	0.00
4) Phenol-d6	8.85	99	399337	10.38	ng	0.00
8) Nitrobenzene-d5	10.85	82	271845	11.39	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	36651	13.91	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	311665	9.11	ng	0.00
22) Terphenyl-d14	22.48	244	415343	9.65	ng	0.00

Target Compounds

						Qvalue
2) n-Nitrosodimethylamine	3.55	42	221628	13.27	ng	# 84
5) Phenol	8.87	94	435341	10.62	ng	90
6) bis(2-Chloroethyl)ether	9.06	93	318407	10.44	ng	# 99
9) Nitrobenzene	10.89	77	309463	11.48	ng	99
10) 2,4-Dimethylphenol	11.82	122	211720	10.41	ng	97
11) 2,4-Dichlorophenol	12.21	162	150596	9.73	ng	97
12) Hexachlorobutadiene	12.89	225	98996	9.05	ng	99
16) 2,4-Dinitrophenol	16.89	184	17342	22.56	ng	93
18) Hexachlorobenzene	19.23	284	102359	9.93	ng	93
19) Pentachlorophenol	19.65	266	39658	20.00	ng	97
21) Benzidine	22.14	184	383031	12.14	ng	99
23) Benzo(a)anthracene	23.83	228	3286058	9.31	ng	98
24) 3,3'-Dichlorobenzidine	23.81	252	348397	12.37	ng	96
25) Chrysene	23.87	228	3743672	8.59	ng	96
26) Indeno(1,2,3-cd)pyrene	27.07	276	1411226	14.00	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	952848	15.68	ng	100
29) Benzo(k)fluoranthene	25.13	252	1223796	8.90	ng	97
30) Benzo(a)pyrene	25.49	252	1095895	14.88	ng	# 95
31) Dibenzo(a,h)anthracene	27.10	278	1143764	15.19	ng	# 95

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

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