

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
 Data File : BE089431.D
 Acq On : 30 Jan 2015 3:36
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 30 07:52:14 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 03:48:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	161714	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	599686	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	251417	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	545560	5.00	ng	0.00
20) Chrysene-d12	23.84	240	698465	5.00	ng	0.00
27) Perylene-d12	25.55	264	744049	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.76	112	56185	0.92	ng	0.02
4) Phenol-d6	8.87	99	68001	0.94	ng	0.03
8) Nitrobenzene-d5	10.85	82	44805	0.92	ng	0.00
14) 2,4,6-Tribromophenol	18.59	330	7486	1.08	ng	0.01
15) 2-Fluorobiphenyl	15.13	172	65141	0.93	ng	0.00
22) Terphenyl-d14	22.48	244	92882	1.15	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.58	42	31391	0.99	ng	# 66
5) Phenol	8.90	94	73902	0.93	ng	93
6) bis(2-Chloroethyl)ether	9.06	93	51921	0.88	ng	# 99
9) Nitrobenzene	10.89	77	52175	0.94	ng	99
10) 2,4-Dimethylphenol	11.84	122	39161	0.99	ng	93
11) 2,4-Dichlorophenol	12.22	162	27847	1.01	ng	90
12) Hexachlorobutadiene	12.89	225	17212	0.93	ng	99
16) 2,4-Dinitrophenol	16.91	184	1618	0.88	ng	# 79
18) Hexachlorobenzene	19.24	284	25240	0.94	ng	93
19) Pentachlorophenol	19.66	266	3437	0.69	ng	94
21) Benzidine	22.15	184	58064	0.95	ng	99
23) Benzo(a)anthracene	23.83	228	638967	1.20	ng	99
24) 3,3'-Dichlorobenzidine	23.82	252	58623	1.41	ng	# 93
25) Chrysene	23.87	228	660407	0.91	ng	94
26) Indeno(1,2,3-cd)pyrene	27.07	276	197968	1.58	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	144666	1.97	ng	98
29) Benzo(k)fluoranthene	25.13	252	198958	1.03	ng	99
30) Benzo(a)pyrene	25.48	252	148995	1.64	ng	99
31) Dibenzo(a,h)anthracene	27.09	278	166745	1.84	ng	97

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

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