

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE022615\
 Data File : BE089715.D
 Acq On : 27 Feb 2015 7:14
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Feb 27 14:23:11 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE022615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 27 13:38:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.33	152	69592	5.00	ng	0.00
7) Naphthalene-d8	12.23	136	302123	5.00	ng	0.00
13) Acenaphthene-d10	16.40	164	140874	5.00	ng	0.00
17) Phenanthrene-d10	19.76	188	256457	5.00	ng	0.00
20) Chrysene-d12	23.65	240	204178	5.00	ng	0.00
27) Perylene-d12	25.34	264	165659	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.49	112	16507	0.98	ng	0.00
4) Phenol-d6	8.61	99	22242	0.97	ng	0.00
8) Nitrobenzene-d5	10.60	82	14276	0.89	ng	0.00
14) 2,4,6-Tribromophenol	18.31	330	1307	0.93	ng	0.00
15) 2-Fluorobiphenyl	14.87	172	29349	0.93	ng	0.00
22) Terphenyl-d14	22.30	244	23970	0.97	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.35	42	9176	0.92	ng	97
5) Phenol	8.64	94	24266	0.97	ng	# 60
6) bis(2-Chloroethyl)ether	8.82	93	18604	0.95	ng	# 99
9) Nitrobenzene	10.64	77	16734	0.94	ng	99
10) 2,4-Dimethylphenol	11.57	122	13484	0.97	ng	99
11) 2,4-Dichlorophenol	11.95	162	9929	0.98	ng	99
12) Hexachlorobutadiene	12.62	225	5256	0.91	ng	100
16) 2,4-Dinitrophenol	16.65	184	214	0.79	ng	93
18) Hexachlorobenzene	18.98	284	6082	0.91	ng	98
19) Pentachlorophenol	19.43	266	962	1.20	ng	98
21) Benzidine	21.96	184	8033	0.92	ng	98
23) Benzo(a)anthracene	23.63	228	140832	0.97	ng	92
24) 3,3'-Dichlorobenzidine	23.63	252	7661	0.98	ng	100
25) Chrysene	23.68	228	161397	0.90	ng	100
26) Indeno(1,2,3-cd)pyrene	26.74	276	23543	1.02	ng	98
28) Benzo(b)fluoranthene	24.91	252	17731	0.96	ng	100
29) Benzo(k)fluoranthene	24.94	252	38679	0.95	ng	99
30) Benzo(a)pyrene	25.27	252	19162	0.96	ng	99
31) Dibenzo(a,h)anthracene	26.77	278	17961	1.01	ng	98

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

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE022615\
Data File : BE089715.D
Acq On : 27 Feb 2015 7:14
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDCCC001EC

Quant Time: Feb 27 14:23:11 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE022615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 27 13:38:44 2015
Response via : Initial Calibration

