

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089394.D
 Acq On : 28 Jan 2015 18:34
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.5

Quant Time: Jan 29 12:10:07 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 29 11:58:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	131394	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	472456	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	192817	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	382156	5.00	ng	0.00
20) Chrysene-d12	23.84	240	427696	5.00	ng	0.00
27) Perylene-d12	25.55	264	437096	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.73	112	19596	0.48	ng	0.00
4) Phenol-d6	8.84	99	24925	0.49	ng	0.00
8) Nitrobenzene-d5	10.84	82	16025	0.48	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	2066	0.01	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	28424	0.01	ng	0.00
22) Terphenyl-d14	22.48	244	25760	0.54	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.59	42	9375	0.40	ng	# 79
5) Phenol	8.87	94	26678	0.50	ng	# 56
6) bis(2-Chloroethyl)ether	9.06	93	20076	0.50	ng	# 99
9) Nitrobenzene	10.88	77	18018	0.48	ng	99
10) 2,4-Dimethylphenol	11.82	122	13848	0.48	ng	97
11) 2,4-Dichlorophenol	12.21	162	10539	0.48	ng	96
12) Hexachlorobutadiene	12.89	225	7536	0.48	ng	100
16) 2,4-Dinitrophenol	16.90	184	269	0.03	ng	# 82
18) Hexachlorobenzene	19.23	284	8596	0.49	ng	98
19) Pentachlorophenol	19.65	266	901	0.27	ng	97
21) Benzidine	22.14	184	15162	0.30	ng	98
23) Benzo(a)anthracene	23.83	228	189121	0.48	ng	99
24) 3,3'-Dichlorobenzidine	23.81	252	11950	0.38	ng	98
25) Chrysene	23.87	228	242323	0.50	ng	100
26) Indeno(1,2,3-cd)pyrene	27.06	276	36896	0.33	ng	# 100
28) Benzo(b)fluoranthene	25.09	252	23241m	0.40	ng	
29) Benzo(k)fluoranthene	25.13	252	58881m	0.42	ng	
30) Benzo(a)pyrene	25.48	252	26303	0.39	ng	99
31) Dibenzo(a,h)anthracene	27.08	278	26099	0.36	ng	97

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

```
Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\  
Data File : BE089394.D  
Acq On    : 28 Jan 2015   18:34  
Operator  : TP/IZ  
Sample    : SSTDICC0.5  
Misc      :  
ALS Vial  : 4      Sample Multiplier: 1
```

Instrument :
BNA_E
LabSampleId :
SSTDICC0.5

Quant Time: Jan 29 12:10:07 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 29 11:58:35 2015
Response via : Initial Calibration

