

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089396.D
 Acq On : 28 Jan 2015 19:52
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICCC001

Quant Time: Jan 29 12:11:43 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	143079	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	509915	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	206344	5.00	ng	0.00
17) Phenanthrene-d10	19.98	188	413648	5.00	ng	0.00
20) Chrysene-d12	23.84	240	509738	5.00	ng	0.00
27) Perylene-d12	25.55	264	539120	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	6.73	112	41287	0.92	ng	0.00
4) Phenol-d6	8.84	99	51342	0.93	ng	0.00
8) Nitrobenzene-d5	10.84	82	33384	0.93	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	4561	0.03	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	58259	0.03	ng	0.00
22) Terphenyl-d14	22.48	244	59923	1.06	ng	0.00

Target Compounds				Qvalue		
2) n-Nitrosodimethylamine	3.57	42	22345	0.87	ng	100
5) Phenol	8.86	94	55468	0.95	ng	100
6) bis(2-Chloroethyl)ether	9.06	93	41188	0.94	ng	# 100
9) Nitrobenzene	10.88	77	38323	0.94	ng	100
10) 2,4-Dimethylphenol	11.82	122	29106	0.94	ng	100
11) 2,4-Dichlorophenol	12.21	162	22551	0.95	ng	100
12) Hexachlorobutadiene	12.89	225	15390	0.91	ng	100
16) 2,4-Dinitrophenol	16.89	184	1078	0.10	ng	100
18) Hexachlorobenzene	19.23	284	17464	0.92	ng	100
19) Pentachlorophenol	19.65	266	2982	0.81	ng	100
21) Benzidine	22.14	184	36001	0.60	ng	100
23) Benzo(a)anthracene	23.83	228	449435	0.95	ng	100
24) 3,3'-Dichlorobenzidine	23.81	252	33656	0.90	ng	100
25) Chrysene	23.87	228	547754	0.95	ng	100
26) Indeno(1,2,3-cd)pyrene	27.06	276	106181	0.79	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	65853m	0.93	ng	
29) Benzo(k)fluoranthene	25.13	252	155556m	0.91	ng	
30) Benzo(a)pyrene	25.48	252	76421	0.92	ng	100
31) Dibenzo(a,h)anthracene	27.08	278	75903	0.86	ng	100

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

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