

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089395.D
 Acq On : 28 Jan 2015 19:13
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.8

Quant Time: Jan 29 12:11:00 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	119094	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	429231	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	178484	5.00	ng	0.00
17) Phenanthrene-d10	19.98	188	360691	5.00	ng	0.00
20) Chrysene-d12	23.84	240	428767	5.00	ng	0.00
27) Perylene-d12	25.55	264	440897	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.73	112	28122	0.76	ng	0.00
4) Phenol-d6	8.84	99	35732	0.78	ng	0.00
8) Nitrobenzene-d5	10.84	82	23028	0.76	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	3208	0.02	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	41387	0.02	ng	0.00
22) Terphenyl-d14	22.48	244	41189	0.87	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.58	42	14472	0.68	ng	# 53
5) Phenol	8.86	94	38251	0.78	ng	100
6) bis(2-Chloroethyl)ether	9.06	93	28965	0.80	ng	# 100
9) Nitrobenzene	10.89	77	26293	0.77	ng	99
10) 2,4-Dimethylphenol	11.82	122	20139	0.77	ng	99
11) 2,4-Dichlorophenol	12.21	162	15604	0.78	ng	98
12) Hexachlorobutadiene	12.89	225	10862	0.76	ng	100
16) 2,4-Dinitrophenol	16.90	184	564	0.06	ng	96
18) Hexachlorobenzene	19.23	284	12746	0.77	ng	95
19) Pentachlorophenol	19.65	266	1796	0.56	ng	99
21) Benzidine	22.14	184	23722	0.47	ng	99
23) Benzo(a)anthracene	23.83	228	304482	0.77	ng	100
24) 3,3'-Dichlorobenzidine	23.81	252	21110	0.67	ng	99
25) Chrysene	23.87	228	383711	0.79	ng	100
26) Indeno(1,2,3-cd)pyrene	27.06	276	65781	0.58	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	39768m	0.69	ng	
29) Benzo(k)fluoranthene	25.13	252	106923m	0.76	ng	
30) Benzo(a)pyrene	25.48	252	47690	0.70	ng	99
31) Dibenzo(a,h)anthracene	27.08	278	46756	0.65	ng	99

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

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