

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021715\
 Data File : BE089601.D
 Acq On : 17 Feb 2015 19:03
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC010

Quant Time: Feb 18 13:44:26 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE021715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 00:56:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	65223	5.00	ng	0.00
7) Naphthalene-d8	12.24	136	262481	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	134173	5.00	ng	0.00
17) Phenanthrene-d10	19.77	188	260298	5.00	ng	0.00
20) Chrysene-d12	23.66	240	300108	5.00	ng	0.00
27) Perylene-d12	25.34	264	248515	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.50	112	176893	9.70	ng	0.00
4) Phenol-d6	8.62	99	234010	9.59	ng	0.00
8) Nitrobenzene-d5	10.61	82	147079	13.59	ng	0.00
14) 2,4,6-Tribromophenol	18.32	330	39950	14.38	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	348435	9.66	ng	0.00
22) Terphenyl-d14	22.30	244	429705	10.57	ng	0.00
Target Compounds						
2) n-Nitrosodimethylamine	3.32	42	85602	9.15	ng	Qvalue # 77
5) Phenol	8.65	94	247535	9.45	ng	85
6) bis(2-Chloroethyl)ether	8.83	93	188980	9.41	ng	# 99
9) Nitrobenzene	10.65	77	175861	13.69	ng	96
10) 2,4-Dimethylphenol	11.58	122	154681	10.45	ng	97
11) 2,4-Dichlorophenol	11.96	162	131615	11.31	ng	99
12) Hexachlorobutadiene	12.63	225	79310	9.85	ng	99
18) Hexachlorobenzene	19.00	284	111624	10.17	ng	90
23) Benzo(a)anthracene	23.64	228	2987138	10.60	ng	97
25) Chrysene	23.69	228	2780943	9.60	ng	93
29) Benzo(k)fluoranthene	24.95	252	784435	11.76	ng	98

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

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