

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089397.D
 Acq On : 28 Jan 2015 20:31
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC002

Quant Time: Jan 29 12:12:45 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	99029	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	340375	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	127882	5.00	ng	0.00
17) Phenanthrene-d10	19.98	188	237950	5.00	ng	0.00
20) Chrysene-d12	23.84	240	344871	5.00	ng	0.00
27) Perylene-d12	25.55	264	390331	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.73	112	63711	2.06	ng	0.00
4) Phenol-d6	8.84	99	77911	2.04	ng	0.00
8) Nitrobenzene-d5	10.84	82	50164	2.09	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	6248	0.06	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	82146	0.06	ng	0.00
22) Terphenyl-d14	22.48	244	80466	2.11	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.56	42	35430	2.00	ng	# 73
5) Phenol	8.86	94	83147	2.05	ng	97
6) bis(2-Chloroethyl)ether	9.06	93	61738	2.05	ng	# 99
9) Nitrobenzene	10.88	77	57313	2.12	ng	99
10) 2,4-Dimethylphenol	11.82	122	43335	2.10	ng	98
11) 2,4-Dichlorophenol	12.21	162	33068	2.08	ng	99
12) Hexachlorobutadiene	12.89	225	23359	2.07	ng	99
16) 2,4-Dinitrophenol	16.89	184	2004	0.30	ng	95
18) Hexachlorobenzene	19.23	284	22740	2.08	ng	99
19) Pentachlorophenol	19.65	266	4959	2.34	ng	97
21) Benzidine	22.14	184	58444	1.44	ng	99
23) Benzo(a)anthracene	23.83	228	678023	2.12	ng	98
24) 3,3'-Dichlorobenzidine	23.81	252	58401	2.30	ng	100
25) Chrysene	23.87	228	816423	2.09	ng	99
26) Indeno(1,2,3-cd)pyrene	27.07	276	206569	2.27	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	124531m	2.43	ng	
29) Benzo(k)fluoranthene	25.13	252	257754m	2.07	ng	
30) Benzo(a)pyrene	25.48	252	143674	2.39	ng	99
31) Dibenzo(a,h)anthracene	27.08	278	150646	2.35	ng	99

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

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