

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020215\
 Data File : BE089457.D
 Acq On : 2 Feb 2015 15:53
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC002

Quant Time: Feb 03 03:15:07 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE020215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 03 03:06:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	36779	5.00	ng	0.00
7) Naphthalene-d8	12.49	136	134405	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	50572	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	94465	5.00	ng	0.00
20) Chrysene-d12	23.84	240	133144	5.00	ng	0.00
27) Perylene-d12	25.56	264	151698	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.76	112	25289	1.98	ng	0.00
4) Phenol-d6	8.87	99	29990	1.94	ng	0.00
8) Nitrobenzene-d5	10.85	82	20400	1.88	ng	0.00
14) 2,4,6-Tribromophenol	18.59	330	2769	1.63	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	28503	2.04	ng	0.00
22) Terphenyl-d14	22.48	244	33092	1.98	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.59	42	13665	2.06	ng	# 88
5) Phenol	8.90	94	32164	1.91	ng	98
6) bis(2-Chloroethyl)ether	9.06	93	22849	1.83	ng	# 97
9) Nitrobenzene	10.89	77	23624	1.90	ng	99
10) 2,4-Dimethylphenol	11.83	122	16848	1.90	ng	99
11) 2,4-Dichlorophenol	12.21	162	12672	1.94	ng	98
12) Hexachlorobutadiene	12.87	225	8478	1.92	ng	98
16) 2,4-Dinitrophenol	16.91	184	732	1.14	ng	# 81
18) Hexachlorobenzene	19.23	284	9959	1.98	ng	100
19) Pentachlorophenol	19.67	266	1136	1.29	ng	96
21) Benzidine	22.15	184	25981	2.84	ng	99
23) Benzo(a)anthracene	23.83	228	235993	2.26	ng	100
24) 3,3'-Dichlorobenzidine	23.82	252	22712	2.93	ng	97
25) Chrysene	23.87	228	260879	1.93	ng	99
26) Indeno(1,2,3-cd)pyrene	27.07	276	88757	4.12	ng	99
28) Benzo(b)fluoranthene	25.10	252	56370	4.13	ng	98
29) Benzo(k)fluoranthene	25.13	252	81413	1.98	ng	99
30) Benzo(a)pyrene	25.49	252	58894	3.19	ng	99
31) Dibenzo(a,h)anthracene	27.10	278	74631	4.13	ng	97

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

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