

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020215\
 Data File : BE089463.D
 Acq On : 3 Feb 2015 1:41
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE020215

Quant Time: Feb 03 11:50:42 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 11:20:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	34184	5.00	ng	0.00
7) Naphthalene-d8	12.49	136	119723	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	48388	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	111573	5.00	ng	0.00
20) Chrysene-d12	23.84	240	155106	5.00	ng	0.00
27) Perylene-d12	25.56	264	171810	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.76	112	11260	0.98	ng	0.00
4) Phenol-d6	8.88	99	12772	0.96	ng	0.00
8) Nitrobenzene-d5	10.85	82	8428	0.97	ng	0.00
14) 2,4,6-Tribromophenol	18.59	330	1501	1.23	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	12411	0.92	ng	0.00
22) Terphenyl-d14	22.48	244	18757	1.07	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.62	42	5697	0.96	ng	# 69
5) Phenol	8.92	94	12786	0.89	ng	# 77
6) bis(2-Chloroethyl)ether	9.06	93	9931	0.95	ng	# 97
9) Nitrobenzene	10.89	77	9831	1.00	ng	99
10) 2,4-Dimethylphenol	11.83	122	7091	1.00	ng	86
11) 2,4-Dichlorophenol	12.22	162	5363	1.00	ng	94
12) Hexachlorobutadiene	12.87	225	3630	0.96	ng	98
16) 2,4-Dinitrophenol	16.92	184	269m	1.02	ng	
18) Hexachlorobenzene	19.23	284	5336	0.91	ng	99
19) Pentachlorophenol	19.67	266	526	0.98	ng	91
21) Benzidine	22.15	184	14179	1.05	ng	98
23) Benzo(a)anthracene	23.83	228	127686	1.03	ng	98
24) 3,3'-Dichlorobenzidine	23.82	252	11181	0.93	ng	98
25) Chrysene	23.87	228	146017	0.97	ng	99
26) Indeno(1,2,3-cd)pyrene	27.07	276	39880	0.80	ng	99
28) Benzo(b)fluoranthene	25.10	252	25355	0.97	ng	99
29) Benzo(k)fluoranthene	25.13	252	45783	1.07	ng	100
30) Benzo(a)pyrene	25.49	252	28697	0.98	ng	99
31) Dibenzo(a,h)anthracene	27.10	278	33674	0.89	ng	99

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

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