

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089400.D
 Acq On : 28 Jan 2015 22:30
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE012815

Quant Time: Jan 29 13:28:21 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 13:20:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	95588	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	327355	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	122334	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	228075	5.00	ng	0.00
20) Chrysene-d12	23.84	240	324089	5.00	ng	0.00
27) Perylene-d12	25.55	264	362067	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.73	112	28275	0.95	ng	0.00
4) Phenol-d6	8.84	99	34519	0.94	ng	0.00
8) Nitrobenzene-d5	10.84	82	22273	0.97	ng	0.00
14) 2,4,6-Tribromophenol	18.57	330	2751	0.99	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	36437	0.98	ng	0.00
22) Terphenyl-d14	22.48	244	33924	0.94	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.58	42	14503	0.92	ng	# 51
5) Phenol	8.86	94	36862	0.94	ng	# 82
6) bis(2-Chloroethyl)ether	9.05	93	27950	0.96	ng	# 100
9) Nitrobenzene	10.89	77	25148	0.97	ng	100
10) 2,4-Dimethylphenol	11.82	122	18699	0.94	ng	97
11) 2,4-Dichlorophenol	12.20	162	14179	0.93	ng	90
12) Hexachlorobutadiene	12.89	225	10442	0.96	ng	99
16) 2,4-Dinitrophenol	16.89	184	538	0.95	ng	88
18) Hexachlorobenzene	19.23	284	10017	0.95	ng	95
19) Pentachlorophenol	19.65	266	1841	0.95	ng	97
21) Benzidine	22.14	184	20804	0.77	ng	100
23) Benzo(a)anthracene	23.83	228	267415	0.89	ng	100
24) 3,3'-Dichlorobenzidine	23.81	252	19714	0.85	ng	99
25) Chrysene	23.87	228	364179	0.99	ng	100
26) Indeno(1,2,3-cd)pyrene	27.06	276	63575	0.87	ng	# 100
28) Benzo(b)fluoranthene	25.09	252	36796m	0.76	ng	
29) Benzo(k)fluoranthene	25.13	252	95494m	0.92	ng	
30) Benzo(a)pyrene	25.48	252	44773	0.82	ng	100
31) Dibenzo(a,h)anthracene	27.08	278	44409	0.81	ng	98

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

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