

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021715\
 Data File : BE089602.D
 Acq On : 17 Feb 2015 19:45
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE021715

Quant Time: Feb 18 14:32:18 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 14:11:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	63854	5.00	ng	0.00
7) Naphthalene-d8	12.24	136	255074	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	127021	5.00	ng	0.00
17) Phenanthrene-d10	19.77	188	251325	5.00	ng	0.00
20) Chrysene-d12	23.66	240	260389	5.00	ng	0.00
27) Perylene-d12	25.34	264	226636	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.50	112	16518	0.99	ng	0.00
4) Phenol-d6	8.62	99	21435	0.98	ng	0.00
8) Nitrobenzene-d5	10.61	82	11301	0.98	ng	0.00
14) 2,4,6-Tribromophenol	18.31	330	2135	0.90	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	32818	0.97	ng	0.00
22) Terphenyl-d14	22.30	244	34683	0.99	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.36	42	7095	0.88	ng	93
5) Phenol	8.65	94	23122	0.98	ng	97
6) bis(2-Chloroethyl)ether	8.83	93	17806	0.95	ng	# 100
9) Nitrobenzene	10.65	77	13951	0.97	ng	99
10) 2,4-Dimethylphenol	11.58	122	13169	1.00	ng	99
11) 2,4-Dichlorophenol	11.96	162	10571	1.00	ng	99
12) Hexachlorobutadiene	12.63	225	7593	0.96	ng	99
16) 2,4-Dinitrophenol	16.65	184	408	1.00	ng	99
18) Hexachlorobenzene	18.99	284	10378	0.96	ng	95
19) Pentachlorophenol	19.44	266	1335	0.93	ng	99
21) Benzidine	21.96	184	11426	0.83	ng	99
23) Benzo(a)anthracene	23.64	228	217894	0.98	ng	100
24) 3,3'-Dichlorobenzidine	23.63	252	11327	1.01	ng	100
25) Chrysene	23.69	228	241262	0.95	ng	99
26) Indeno(1,2,3-cd)pyrene	26.75	276	41944	0.89	ng	# 100
28) Benzo(b)fluoranthene	24.92	252	33276	0.88	ng	100
29) Benzo(k)fluoranthene	24.94	252	61658	0.93	ng	100
30) Benzo(a)pyrene	25.28	252	32790	0.88	ng	99
31) Dibenzo(a,h)anthracene	26.78	278	32385	0.89	ng	100

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

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