

Data Path : Z:\HPCHEM1\BNA E\DATA\BE093015\
 Data File : BE090979.D
 Acq On : 30 Sep 2015 15:17
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Oct 01 03:19:33 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 30 16:01:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	26676	5.00	ng	-0.01
7) Naphthalene-d8	10.23	136	119059	5.00	ng	-0.02
13) Acenaphthene-d10	14.11	164	67069	5.00	ng	0.00
19) Phenanthrene-d10	16.86	188	177868	5.00	ng	0.00
26) Chrysene-d12	21.02	240	226129	5.00	ng	0.00
35) Perylene-d12	23.27	264	198777	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.11	112	4954	0.98	ng	-0.01
5) Phenol-d6	6.68	99	6254	0.92	ng	-0.03
8) Nitrobenzene-d5	8.64	82	6318	0.96	ng	-0.01
14) 2,4,6-Tribromophenol	15.62	330	1953	0.99	ng	-0.02
15) 2-Fluorobiphenyl	12.73	172	16497	0.91	ng	-0.01
29) Terphenyl-d14	19.49	244	24193	0.89	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.08	88	11887	3.39	ng	97
3) n-Nitrosodimethylamine	3.37	42	4772	1.11	ng	98
6) bis(2-Chloroethyl)ether	6.92	93	7430	1.06	ng	97
9) Nitrobenzene	8.68	77	6894	1.01	ng	97
10) Naphthalene	10.28	128	24195	0.99	ng	99
11) Hexachlorobutadiene	10.57	225	3989	0.97	ng	98
12) 2-Methylnaphthalene	11.92	142	14309	0.95	ng	99
16) Acenaphthylene	13.82	152	101271	0.90	ng	100
17) Acenaphthene	14.17	154	16988	0.95	ng	96
18) Fluorene	15.17	166	22317	1.01	ng	99
20) 4-Bromophenyl-phenylether	16.06	248	6068	0.90	ng	87
21) Hexachlorobenzene	16.18	284	30657	0.92	ng	99
22) Pentachlorophenol	16.56	266	1925	0.99	ng	98
23) Phenanthrene	16.89	178	40993	0.96	ng	100
24) Anthracene	17.00	178	34098	0.89	ng	99
25) Fluoranthene	18.91	202	44566	0.93	ng	100
27) Benzidine	19.14	184	3414	1.04	ng	94
28) Pyrene	19.27	202	46907	0.89	ng	100
30) Benzo(a)anthracene	21.00	228	44626	0.89	ng	98
32) Chrysene	21.05	228	57957	0.97	ng	99
33) Bis(2-ethylhexyl)phthalate	20.95	149	13555	0.84	ng	99
34) Indeno(1,2,3-cd)pyrene	25.60	276	46049	0.91	ng	100
36) Benzo(b)fluoranthene	22.59	252	39669	0.91	ng	99
37) Benzo(k)fluoranthene	22.63	252	50273	0.88	ng	100
38) Benzo(a)pyrene	23.17	252	35753	0.88	ng	99
39) Dibenzo(a,h)anthracene	25.62	278	34761	0.87	ng	99
40) Benzo(g,h,i)perylene	26.31	276	39045	0.90	ng	100

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

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