

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020516\
 Data File : BE091395.D
 Acq On : 5 Feb 2016 11:15
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Feb 05 16:38:17 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 00:27:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	44997	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	218478	5.00	ng	0.00
8) Acenaphthene-d10	13.35	164	142107	5.00	ng	0.00
14) Phenanthrene-d10	16.09	188	360681	5.00	ng	0.00
18) Chrysene-d12	20.21	240	408987	5.00	ng	0.00
24) Perylene-d12	22.10	264	387908	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.95	112	7810	0.93	ng	0.00
3) Phenol-d6	6.56	99	12237	1.00	ng	0.00
5) Nitrobenzene-d5	8.17	82	10193	0.94	ng	0.00
9) 2,4,6-Tribromophenol	14.91	330	3840	0.87	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	34519	0.93	ng	0.00
20) Terphenyl-d14	18.62	244	59357	0.98	ng	0.00

Target Compounds				Qvalue		
6) Naphthalene	9.59	128	44934	0.91	ng	100
7) 2-Methylnaphthalene	11.13	142	26075	0.96	ng	96
11) Acenaphthylene	13.10	152	190724	0.92	ng	100
12) Acenaphthene	13.42	154	34773	0.89	ng	99
13) Fluorene	14.40	166	46950	0.93	ng	100
15) Phenanthrene	16.12	178	85560	0.90	ng	99
16) Anthracene	16.21	178	74454	0.94	ng	99
17) Fluoranthene	18.09	202	80733	0.91	ng	99
19) Pyrene	18.47	202	85795	0.97	ng	100
21) Benzo(a)anthracene	20.19	228	69663	0.91	ng	99
22) Chrysene	20.25	228	102336	0.95	ng	100
23) Indeno(1,2,3-cd)pyrene	23.70	276	71786	0.91	ng	100
25) Benzo(b)fluoranthene	21.55	252	77476	0.93	ng	98
26) Benzo(k)fluoranthene	21.57	252	87521	0.92	ng	99
27) Benzo(a)pyrene	22.00	252	68026	0.91	ng	99
28) Dibenzo(a,h)anthracene	23.65	278	59502	0.89	ng	99
29) Benzo(g,h,i)perylene	24.25	276	77837	0.93	ng	99

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

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