

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021716\
 Data File : BE091404.D
 Acq On : 17 Feb 2016 15:01
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Feb 17 23:58:58 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	44281	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	217555	5.00	ng	0.00
8) Acenaphthene-d10	13.35	164	143696	5.00	ng	0.00
14) Phenanthrene-d10	16.08	188	359334	5.00	ng	-0.01
18) Chrysene-d12	20.21	240	396876	5.00	ng	0.00
24) Perylene-d12	22.09	264	384031	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.94	112	7907	0.95	ng	0.00
3) Phenol-d6	6.56	99	12262	1.01	ng	-0.01
5) Nitrobenzene-d5	8.16	82	11418	1.03	ng	0.00
9) 2,4,6-Tribromophenol	14.90	330	4235	0.92	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	36337	0.97	ng	0.00
20) Terphenyl-d14	18.61	244	61276	1.04	ng	0.00

Target Compounds						Qvalue
6) Naphthalene	9.59	128	46227	0.94	ng	99
7) 2-Methylnaphthalene	11.12	142	27495	1.01	ng	99
11) Acenaphthylene	13.10	152	206270	0.97	ng	99
12) Acenaphthene	13.42	154	36533	0.93	ng	98
13) Fluorene	14.40	166	49625	0.97	ng	99
15) Phenanthrene	16.12	178	89486	0.95	ng	99
16) Anthracene	16.21	178	76435	0.97	ng	100
17) Fluoranthene	18.09	202	86580	0.97	ng	99
19) Pyrene	18.47	202	90440	1.05	ng	100
21) Benzo(a)anthracene	20.18	228	69419	0.93	ng	99
22) Chrysene	20.24	228	103220	0.99	ng	99
23) Indeno(1,2,3-cd)pyrene	23.69	276	64263	0.85	ng	99
25) Benzo(b)fluoranthene	21.54	252	79799	0.96	ng	99
26) Benzo(k)fluoranthene	21.57	252	89031	0.94	ng	99
27) Benzo(a)pyrene	22.00	252	69520	0.93	ng	99
28) Dibenzo(a,h)anthracene	23.65	278	53190	0.83	ng	98
29) Benzo(a,h,i)perylene	24.25	276	74956	0.91	ng	99

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

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