

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021716\
 Data File : BE091411.D
 Acq On : 17 Feb 2016 19:20
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Feb 18 00:00:41 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	6.86	152	43942	5.00	ng	-0.01
4) Naphthalene-d8	9.53	136	218633	5.00	ng	-0.01
8) Acenaphthene-d10	13.34	164	138868	5.00	ng	0.00
14) Phenanthrene-d10	16.08	188	331652	5.00	ng	-0.01
18) Chrysene-d12	20.22	240	393856	5.00	ng	0.00
24) Perylene-d12	22.09	264	411540	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	4.91	112	9778	1.13	ng	-0.04
3) Phenol-d6	6.54	99	15154	1.21	ng	-0.03
5) Nitrobenzene-d5	8.14	82	14535	1.23	ng	-0.03
9) 2,4,6-Tribromophenol	14.89	330	5990	1.18	ng	-0.01
10) 2-Fluorobiphenyl	11.99	172	35001	0.97	ng	-0.02
20) Terphenyl-d14	18.61	244	63454	1.09	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	9.57	128	46230	0.94	ng	100
7) 2-Methylnaphthalene	11.10	142	28643	1.04	ng	98
11) Acenaphthylene	13.09	152	229706	1.09	ng	100
12) Acenaphthene	13.41	154	35125	0.92	ng	97
13) Fluorene	14.39	166	47457	0.96	ng	99
15) Phenanthrene	16.12	178	80658	0.93	ng	99
16) Anthracene	16.20	178	79665	1.09	ng	100
17) Fluoranthene	18.08	202	91101	1.09	ng	99
19) Pyrene	18.47	202	93339	1.09	ng	99
21) Benzo(a)anthracene	20.18	228	92947	1.19	ng	99
22) Chrysene	20.24	228	95976	0.93	ng	98
23) Indeno(1,2,3-cd)pyrene	23.69	276	98929	1.23	ng	97
25) Benzo(b)fluoranthene	21.54	252	92406	1.02	ng	98
26) Benzo(k)fluoranthene	21.57	252	104477	1.03	ng	99
27) Benzo(a)pyrene	21.99	252	93144	1.12	ng	99
28) Dibenzo(a,h)anthracene	23.64	278	96129	1.22	ng	99
29) Benzo(a,h,i)perylene	24.23	276	100554	1.10	ng	98

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

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