

Data Path : Z:\HPCHEM1\BNA E\Data\BE101415\
 Data File : BE091011.D
 Acq On : 14 Oct 2015 19:22
 Operator : UM/IZ
 Sample : SSTD1.677
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.677

Quant Time: Oct 15 05:35:35 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE101415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 05:25:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	5980	0.40	ng/ul	0.00
2) Naphthalene-d8	9.62	136	28971	0.40	ng/ul	0.00
6) Acenaphthene-d10	13.44	164	16257	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.16	188	40282	0.40	ng/ul	0.00
16) Chrysene-d12	20.31	240	52997	0.40	ng/ul	0.00
20) Perylene-d12	22.22	264	53413	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.13	152	49216	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.15	212	140794	1.31	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	9.66	128	92944	1.31	ng/ul	100
5) 2-Methylnaphthalene	11.20	142	57539	1.27	ng/ul	96
7) Acenaphthylene	13.18	152	91869	0.35	ng/ul	99
8) Acenaphthene	13.50	153	65098	0.35	ng/ul	98
9) Fluorene	14.49	166	79733	1.22	ng/ul	99
11) Pentachlorophenol	15.92	266	7066	1.34	ng/ul	96
12) Phenanthrene	16.20	178	545248	3.35	ng/ul	98
13) Anthracene	16.28	178	531740	3.39	ng/ul	99
15) Fluoranthene	18.18	202	673860	3.41	ng/ul	100
17) Pyrene	18.56	202	682774	3.33	ng/ul	99
18) Benzo(a)anthracene	20.28	228	177415	1.26	ng/ul	99
19) Chrysene	20.34	228	184979	1.24	ng/ul	98
21) Benzo(b)fluoranthene	21.65	252	195745	1.22	ng/ul	83
22) Benzo(k)fluoranthene	21.68	252	187643	1.21	ng/ul#	84
23) Benzo(a)pyrene	22.12	252	188464	1.26	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	23.87	276	746005	3.25	ng/ul	100
25) Dibenzo(a,h)anthracene	23.83	278	192174	1.22	ng/ul	95
26) Benzo(g,h,i)perylene	24.44	276	798540	3.29	ng/ul#	90

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

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