

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE112817\
 Data File : BE094886.D
 Acq On : 28 Nov 2017 16:26
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.401

Manual Integrations
 APPROVED

Sohil
 11/30/2017 8:47:21 PM

Quant Time: Nov 29 02:01:44 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 29 01:10:42 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	620	0.40	ng/ul	0.01
2) Naphthalene-d8	10.72	136	4308	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3077	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7404	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7562	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7284	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	2854	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	8318	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.77	128	3607	0.352	ng/ul#	95
5) 2-Methylnaphthalene	12.39	142	2768	0.388	ng/ul	98
7) Acenaphthylene	14.25	152	4709	0.334	ng/ul	98
8) Acenaphthene	14.58	153	3471	0.350	ng/ul	99
9) Fluorene	15.58	166	4139	0.344	ng/ul#	94
11) Pentachlorophenol	17.08	266	92m	0.236	ng/ul	
12) Phenanthrene	17.31	178	6927	0.348	ng/ul#	93
13) Anthracene	17.41	178	6842	0.338	ng/ul#	93
15) Fluoranthene	19.32	202	8274	0.317	ng/ul	84
17) Pyrene	19.68	202	8659	0.366	ng/ul#	78
18) Benzo(a)anthracene	21.41	228	8229	0.358	ng/ul#	87
19) Chrysene	21.47	228	8477	0.369	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	7842	0.361	ng/ul	84
22) Benzo(k)fluoranthene	23.24	252	8239	0.357	ng/ul#	87
23) Benzo(a)pyrene	23.86	252	7849	0.364	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.69	276	8653	0.368	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.67	278	7195	0.363	ng/ul#	89
26) Benzo(g,h,i)perylene	27.53	276	7299	0.370	ng/ul	94

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

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