

Data Path : Z:\HPCHEM1\BNA_E\Data\BE121817\
 Data File : BE095005.D
 Acq On : 18 Dec 2017 12:17
 Operator : SJ/JU
 Sample : SST0.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.202

Quant Time: Dec 18 15:23:11 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:20:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4892	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14901	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8027	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	20388	0.40	ng/ul	0.00
16) Chrysene-d12	21.85	240	18894	0.40	ng/ul	0.00
20) Perylene-d12	24.51	264	15965	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	4767	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	11710	0.17	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	7778	0.219	ng/ul#	90
5) 2-Methylnaphthalene	12.92	142	6866	0.278	ng/ul	100
7) Acenaphthylene	14.78	152	7037	0.191	ng/ul	97
8) Acenaphthene	15.11	153	5513	0.213	ng/ul	92
11) Pentachlorophenol	17.41	266	319	0.297	ng/ul#	61
12) Phenanthrene	17.78	178	10038	0.183	ng/ul#	82
13) Anthracene	17.87	178	10162	0.182	ng/ul#	86
15) Fluoranthene	19.76	202	12519	0.174	ng/ul	84
17) Pyrene	20.11	202	13600	0.230	ng/ul#	80
18) Benzo(a)anthracene	21.82	228	10561	0.184	ng/ul#	85
19) Chrysene	21.88	228	11367	0.198	ng/ul	96
21) Benzo(b)fluoranthene	23.68	252	10707	0.225	ng/ul	88
22) Benzo(k)fluoranthene	23.73	252	9949	0.197	ng/ul#	91
23) Benzo(a)pyrene	24.38	252	9128	0.193	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	27.34	276	10036	0.195	ng/ul#	82
25) Dibenzo(a,h)anthracene	27.37	278	8059	0.185	ng/ul	93
26) Benzo(g,h,i)perylene	28.22	276	8697	0.201	ng/ul	97

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

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