

Data Path : U:\HPCHEM1\BNA E\DATA\BE011918\
 Data File : BE095076.D
 Acq On : 19 Jan 2018 15:14
 Operator : SJ/JU
 Sample : SSTD1.605
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.605

Quant Time: Jan 19 18:03:20 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE011918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 19 17:59:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4851	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	15367	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8884	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	19233	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	21204	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	16170	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	41228	1.62	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	105383	1.72	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.34	128	64068	1.574	ng/ul#	88
5) 2-Methylnaphthalene	12.91	142	46433	1.627	ng/ul	99
7) Acenaphthylene	14.77	152	65741	1.621	ng/ul#	96
8) Acenaphthene	15.11	153	49082	1.578	ng/ul	95
9) Fluorene	16.07	166	53149	1.598	ng/ul	87
11) Pentachlorophenol	17.39	266	1853	1.603	ng/ul#	56
12) Phenanthrene	17.78	178	87036	1.684	ng/ul#	83
13) Anthracene	17.87	178	84745	1.833	ng/ul#	87
15) Fluoranthene	19.76	202	117259	1.748	ng/ul	84
17) Pyrene	20.11	202	116414	1.523	ng/ul#	80
18) Benzo(a)anthracene	21.82	228	96380	1.617	ng/ul#	86
19) Chrysene	21.88	228	101873	1.562	ng/ul	95
21) Benzo(b)fluoranthene	23.67	252	89529	1.634	ng/ul	82
22) Benzo(k)fluoranthene	23.72	252	91164	1.768	ng/ul#	84
23) Benzo(a)pyrene	24.37	252	81348	1.711	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.33	276	79064	1.824	ng/ul#	81
25) Dibenzo(a,h)anthracene	27.36	278	60897	1.923	ng/ul#	87
26) Benzo(g,h,i)perylene	28.21	276	72579	1.726	ng/ul	93

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

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