

Data Path : U:\HPCHEM1\BNA E\DATA\BE011918\
 Data File : BE095072.D
 Acq On : 19 Jan 2018 12:51
 Operator : SJ/JU
 Sample : SST00.101
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.101

Quant Time: Jan 19 15:25:09 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 15:16:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4563	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14512	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	7770	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	18324	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	15716	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	13345	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	2326	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	5417	0.09	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.34	128	3850	0.100	ng/ul#	91
5) 2-Methylnaphthalene	12.91	142	2555	0.086	ng/ul	97
7) Acenaphthylene	14.78	152	3360	0.095	ng/ul	99
8) Acenaphthene	15.11	153	2708	0.099	ng/ul	95
9) Fluorene	16.07	166	2743	0.090	ng/ul	91
12) Phenanthrene	17.78	178	4832	0.100	ng/ul#	82
13) Anthracene	17.87	178	3845	0.084	ng/ul#	87
15) Fluoranthene	19.76	202	5753	0.091	ng/ul	84
17) Pyrene	20.11	202	5992	0.110	ng/ul#	83
18) Benzo(a)anthracene	21.82	228	4301	0.095	ng/ul#	85
19) Chrysene	21.89	228	4870	0.102	ng/ul	98
21) Benzo(b)fluoranthene	23.68	252	4057	0.087	ng/ul	95
22) Benzo(k)fluoranthene	23.73	252	3643	0.086	ng/ul	97
23) Benzo(a)pyrene	24.38	252	3474	0.087	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	27.33	276	2731	0.062	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.37	278	1751	0.049	ng/ul#	88
26) Benzo(a,h,i)perylene	28.23	276	3059	0.080	ng/ul	96

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

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