

Data Path : U:\HPCHEM1\BNA E\DATA\BE011918\
 Data File : BE095073.D
 Acq On : 19 Jan 2018 13:26
 Operator : SJ/JU
 Sample : SST00.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.202

Quant Time: Jan 19 18:10:09 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 18:05:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4813	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	15406	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8411	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	19252	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	16507	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	13656	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	4937	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	11236	0.18	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.34	128	7989	0.196	ng/ul#	87
5) 2-Methylnaphthalene	12.91	142	5494	0.192	ng/ul	100
7) Acenaphthylene	14.77	152	7252	0.189	ng/ul	96
8) Acenaphthene	15.11	153	5695	0.193	ng/ul	92
9) Fluorene	16.07	166	5939	0.189	ng/ul	92
11) Pentachlorophenol	17.41	266	152	0.138	ng/ul#	81
12) Phenanthrene	17.78	178	9889	0.191	ng/ul#	81
13) Anthracene	17.87	178	7997	0.173	ng/ul#	86
15) Fluoranthene	19.76	202	12062	0.180	ng/ul	85
17) Pyrene	20.11	202	12318	0.207	ng/ul#	80
18) Benzo(a)anthracene	21.82	228	8824	0.190	ng/ul#	84
19) Chrysene	21.89	228	10177	0.200	ng/ul	98
21) Benzo(b)fluoranthene	23.68	252	8914	0.193	ng/ul	88
22) Benzo(k)fluoranthene	23.73	252	7950	0.183	ng/ul#	91
23) Benzo(a)pyrene	24.38	252	7377	0.184	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	27.34	276	6213	0.170	ng/ul#	82
25) Dibenzo(a,h)anthracene	27.37	278	4242	0.159	ng/ul	98
26) Benzo(g,h,i)perylene	28.22	276	6409	0.180	ng/ul	96

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

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