

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033020\
 Data File : BM025854.D
 Acq On : 30 Mar 2020 11:24
 Operator : CG/JU
 Sample : SST0.281
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.281

Quant Time: Mar 30 12:34:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM033020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 30 12:31:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1548	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	5803	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3498	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	7760	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	6772	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	6974	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	2093	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	4689	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	3660	0.210	ng/ul	98
5) 2-Methylnaphthalene	12.01	142	2460	0.193	ng/ul	100
7) Acenaphthylene	13.92	152	3067	0.187	ng/ul	98
8) Acenaphthene	14.27	153	2604	0.205	ng/ul	98
9) Fluorene	15.26	166	3022	0.190	ng/ul	98
11) Pentachlorophenol	16.62	266	291	0.143	ng/ul	94
12) Phenanthrene	16.99	178	5114	0.204	ng/ul	98
13) Anthracene	17.09	178	4113	0.185	ng/ul	96
15) Fluoranthene	19.01	202	5726	0.183	ng/ul	97
17) Pyrene	19.37	202	5789	0.225	ng/ul	97
18) Benzo(a)anthracene	21.13	228	4647	0.187	ng/ul	98
19) Chrysene	21.18	228	5830	0.213	ng/ul	98
21) Benzo(b)fluoranthene	22.66	252	5472	0.196	ng/ul	92
22) Benzo(k)fluoranthene	22.70	252	5810	0.199	ng/ul#	94
23) Benzo(a)pyrene	23.20	252	4604	0.197	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.42	276	5635	0.193	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.43	278	4378	0.181	ng/ul	93
26) Benzo(g,h,i)perylene	26.06	276	5608	0.226	ng/ul	98

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

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