

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM033020\
 Data File : BM025858.D
 Acq On : 30 Mar 2020 13:48
 Operator : CG/JU
 Sample : SSTD3.279
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.279

Quant Time: Mar 30 14:19:59 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	1630	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	6032	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3794	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8390	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	9398	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	9001	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	34074	3.02	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	86509	3.17	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.38	128	57635	3.175	ng/ul	97
5) 2-Methylnaphthalene	12.00	142	40198	3.040	ng/ul	100
7) Acenaphthylene	13.92	152	53769	3.026	ng/ul	97
8) Acenaphthene	14.27	153	43375	3.148	ng/ul	98
9) Fluorene	15.26	166	53052	3.071	ng/ul	99
11) Pentachlorophenol	16.61	266	6936	3.158	ng/ul	94
12) Phenanthrene	16.99	178	88086	3.256	ng/ul	99
13) Anthracene	17.08	178	81272	3.385	ng/ul	98
15) Fluoranthene	19.01	202	111077	3.280	ng/ul	98
17) Pyrene	19.37	202	108269	3.038	ng/ul	99
18) Benzo(a)anthracene	21.13	228	107365	3.108	ng/ul	99
19) Chrysene	21.18	228	113602	2.993	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	115195	3.199	ng/ul	93
22) Benzo(k)fluoranthene	22.70	252	118336	3.133	ng/ul#	93
23) Benzo(a)pyrene	23.20	252	98483	3.257	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.41	276	129802	3.444	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.42	278	103766	3.323	ng/ul	92
26) Benzo(g,h,i)perylene	26.05	276	116600	3.643	ng/ul	97

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

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