

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM040220\
 Data File : BM025891.D
 Acq On : 01 Apr 2020 17:22
 Operator : CG/JU
 Sample : SSTD3.280
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.280

Quant Time: Apr 01 17:53:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 16:04:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2362	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9905	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	6835	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15739	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17545	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	17654	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	58107	3.43	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	168026	3.67	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	92849	3.153	ng/ul	98
5) 2-Methylnaphthalene	12.00	142	68011	3.434	ng/ul	99
7) Acenaphthylene	13.92	152	100638	3.470	ng/ul	98
8) Acenaphthene	14.27	153	79029	3.261	ng/ul	99
9) Fluorene	15.26	166	99328	3.464	ng/ul	98
11) Pentachlorophenol	16.61	266	15527	4.767	ng/ul	96
12) Phenanthrene	16.99	178	165374	3.351	ng/ul	98
13) Anthracene	17.08	178	157725	3.821	ng/ul	98
15) Fluoranthene	19.01	202	210314	3.675	ng/ul	98
17) Pyrene	19.37	202	207315	3.062	ng/ul	99
18) Benzo(a)anthracene	21.13	228	214969	3.727	ng/ul	99
19) Chrysene	21.18	228	216831	3.116	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	224805	3.290	ng/ul	97
22) Benzo(k)fluoranthene	22.69	252	231758	3.312	ng/ul	96
23) Benzo(a)pyrene	23.20	252	200517	3.567	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.41	276	260368	3.643	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.42	278	217540	3.934	ng/ul	98
26) Benzo(g,h,i)perylene	26.05	276	215652	3.101	ng/ul	100

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

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