

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082820\
 Data File : BM027284.D
 Acq On : 28 Aug 2020 09:36
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Aug 28 11:51:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 10:39:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2181	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7243	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	4578	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	9940	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	7266	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	6208	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5130	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	10498	0.37	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.41	128	8634	0.398	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	6049	0.407	ng/ul	96
7) Acenaphthylene	13.95	152	8026	0.404	ng/ul	98
8) Acenaphthene	14.29	153	6977	0.404	ng/ul	99
9) Fluorene	15.29	166	8249	0.399	ng/ul	99
11) Pentachlorophenol	16.65	266	740	0.226	ng/ul	94
12) Phenanthrene	17.02	178	13219	0.393	ng/ul	98
13) Anthracene	17.11	178	11607	0.380	ng/ul	99
15) Fluoranthene	19.05	202	14611	0.362	ng/ul	99
17) Pyrene	19.41	202	14448	0.465	ng/ul	99
18) Benzo(a)anthracene	21.17	228	10935	0.379	ng/ul	100
19) Chrysene	21.22	228	12717	0.407	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	10868	0.408	ng/ul	99
22) Benzo(k)fluoranthene	22.76	252	11738	0.424	ng/ul	99
23) Benzo(a)pyrene	23.27	252	9600	0.398	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.54	276	12759	0.384	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.55	278	10486	0.387	ng/ul	99
26) Benzo(g,h,i)perylene	26.19	276	11411	0.398	ng/ul	98

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

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