

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082020\
 Data File : BM027145.D
 Acq On : 20 Aug 2020 01:44
 Operator : JU/CG
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.430

Quant Time: Aug 20 06:54:56 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 : Wed Aug 19 11:15:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1378	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	4304	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2609	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	5728	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	5262	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	5021	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3077	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	6683	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	5410	0.420	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	3671	0.416	ng/ul	100
7) Acenaphthylene	13.98	152	4552	0.402	ng/ul	99
8) Acenaphthene	14.32	153	4201	0.427	ng/ul	99
9) Fluorene	15.31	166	4954	0.420	ng/ul	100
11) Pentachlorophenol	16.67	266	699	0.370	ng/ul	94
12) Phenanthrene	17.05	178	8252	0.426	ng/ul	100
13) Anthracene	17.14	178	7160	0.407	ng/ul	99
15) Fluoranthene	19.07	202	9632	0.414	ng/ul	99
17) Pyrene	19.43	202	9760	0.434	ng/ul	99
18) Benzo(a)anthracene	21.18	228	8481	0.406	ng/ul	99
19) Chrysene	21.23	228	9950	0.440	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	9340	0.434	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	10284	0.459	ng/ul	99
23) Benzo(a)pyrene	23.29	252	8463	0.434	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.57	276	10936	0.407	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.58	278	8908	0.407	ng/ul	98
26) Benzo(g,h,i)perylene	26.23	276	9148	0.394	ng/ul	99

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

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