

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081720\
 Data File : BM027111.D
 Acq On : 17 Aug 2020 11:53
 Operator : JU/CG
 Sample : SSTD0.477
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.477

Quant Time: Aug 17 12:40:26 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 : Mon Aug 17 12:32:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1314	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	4001	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2456	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5396	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	5199	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	5337	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2870	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	6559	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	5101	0.405	ng/ul	100
5) 2-Methylnaphthalene	12.07	142	3426	0.411	ng/ul	100
7) Acenaphthylene	13.98	152	4599	0.402	ng/ul	100
8) Acenaphthene	14.33	153	3923	0.405	ng/ul	100
9) Fluorene	15.32	166	4714	0.416	ng/ul	100
11) Pentachlorophenol	16.67	266	767	0.563	ng/ul	100
12) Phenanthrene	17.05	178	7806	0.424	ng/ul	100
13) Anthracene	17.14	178	7063	0.456	ng/ul	100
15) Fluoranthene	19.07	202	9313	0.435	ng/ul	100
17) Pyrene	19.43	202	9477	0.378	ng/ul	100
18) Benzo(a)anthracene	21.19	228	8676	0.441	ng/ul	100
19) Chrysene	21.24	228	9532	0.390	ng/ul	100
21) Benzo(b)fluoranthene	22.74	252	9838	0.404	ng/ul	100
22) Benzo(k)fluoranthene	22.79	252	10150	0.379	ng/ul	100
23) Benzo(a)pyrene	23.30	252	8703	0.426	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.58	276	12258	0.545	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	10070	0.582	ng/ul	100
26) Benzo(g,h,i)perylene	26.25	276	10691	0.519	ng/ul	100

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

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