

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072020\
 Data File : BN011203.D
 Acq On : 20 Jul 2020 10:34
 Operator : CG/JU
 Sample : SSTDO.401
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.401

Quant Time: Jul 20 11:27:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 20 11:25:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	1138	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	3908	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	2275	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	4634	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	3040	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	2742	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	4105	0.45	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	8164	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.29	128	4959	0.430	ng/ul	100
5) 2-Methylnaphthalene	11.92	142	3262	0.451	ng/ul	100
7) Acenaphthylene	13.84	152	4632	0.457	ng/ul	100
8) Acenaphthene	14.19	153	3668	0.432	ng/ul	100
9) Fluorene	15.18	166	4085	0.437	ng/ul	100
11) Pentachlorophenol	16.55	266	394	0.363	ng/ul	99
12) Phenanthrene	16.92	178	6384	0.423	ng/ul	100
13) Anthracene	17.01	178	5223	0.419	ng/ul	100
15) Fluoranthene	18.95	202	7004	0.414	ng/ul	100
17) Pyrene	19.31	202	6832	0.517	ng/ul	100
18) Benzo(a)anthracene	21.08	228	4483	0.415	ng/ul	100
19) Chrysene	21.13	228	5990	0.436	ng/ul	100
21) Benzo(b)fluoranthene	22.60	252	4908	0.402	ng/ul	100
22) Benzo(k)fluoranthene	22.64	252	5727	0.443	ng/ul	100
23) Benzo(a)pyrene	23.14	252	4266	0.414	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.32	276	5101	0.361	ng/ul	99
25) Dibenzo(a,h)anthracene	25.34	278	3859	0.334	ng/ul	100
26) Benzo(g,h,i)perylene	25.94	276	5000	0.382	ng/ul	100

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

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