

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120118\
 Data File : BN003653.D
 Acq On : 01 Dec 2018 09:56
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.477

Quant Time: Dec 03 03:32:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 03 03:30:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5115	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	19336	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	9934	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	22670	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	20701	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	18380	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	11561	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	27445	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	22306	0.411	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	15390	0.406	ng/ul	100
7) Acenaphthylene	14.19	152	18723	0.419	ng/ul	100
8) Acenaphthene	14.54	153	16347	0.417	ng/ul	100
9) Fluorene	15.52	166	18400	0.403	ng/ul	99
11) Pentachlorophenol	16.86	266	1500	0.367	ng/ul	99
12) Phenanthrene	17.26	178	31428	0.422	ng/ul	99
13) Anthracene	17.35	178	25199	0.412	ng/ul	100
15) Fluoranthene	19.27	202	38114	0.434	ng/ul	97
17) Pyrene	19.64	202	36502	0.492	ng/ul	100
18) Benzo(a)anthracene	21.38	228	27412	0.419	ng/ul	98
19) Chrysene	21.43	228	30620	0.409	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	31247	0.411	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	29490	0.398	ng/ul	99
23) Benzo(a)pyrene	23.60	252	26351	0.400	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	27387	0.363	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.07	278	20140	0.331	ng/ul	98
26) Benzo(g,h,i)perylene	26.78	276	24771	0.377	ng/ul	98

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

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