

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012785.D
 Acq On : 01 Dec 2020 17:03
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Dec 01 17:39:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 01 17:05:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	829	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	3360	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1973	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3952	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3890	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	4017	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	2082	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	4122	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.24	128	4226	0.418	ng/ul	98
5) 2-Methylnaphthalene	11.86	142	2909	0.450	ng/ul	100
7) Acenaphthylene	13.80	152	4162	0.433	ng/ul	98
8) Acenaphthene	14.14	153	3184	0.393	ng/ul	99
9) Fluorene	15.14	166	3518	0.395	ng/ul	100
11) Pentachlorophenol	16.49	266	599	0.477	ng/ul	96
12) Phenanthrene	16.88	178	5482	0.399	ng/ul	98
13) Anthracene	16.97	178	5028	0.421	ng/ul	98
15) Fluoranthene	18.91	202	6972	0.439	ng/ul	99
17) Pyrene	19.27	202	7293	0.419	ng/ul	97
18) Benzo(a)anthracene	21.04	228	7432	0.533	ng/ul	99
19) Chrysene	21.09	228	8278	0.475	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	8321	0.450	ng/ul	96
22) Benzo(k)fluoranthene	22.58	252	8947	0.420	ng/ul	96
23) Benzo(a)pyrene	23.07	252	7838	0.451	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.24	276	9936	0.437	ng/ul	99
25) Dibenzo(a,h)anthracene	25.25	278	7965	0.436	ng/ul	95
26) Benzo(g,h,i)perylene	25.87	276	8343	0.408	ng/ul	98

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

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