

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012943.D
 Acq On : 07 Dec 2020 17:30
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.429

Quant Time: Dec 07 18:00:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 12:24:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	859	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	3523	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1985	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4018	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	3204	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3043	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	2093	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	4266	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	4024	0.395	ng/ul	94
5) 2-Methylnaphthalene	11.84	142	2703	0.404	ng/ul	98
7) Acenaphthylene	13.77	152	3854	0.410	ng/ul	95
8) Acenaphthene	14.12	153	3076	0.399	ng/ul	98
9) Fluorene	15.12	166	3385	0.383	ng/ul	98
11) Pentachlorophenol	16.47	266	472	0.406	ng/ul	95
12) Phenanthrene	16.85	178	5407	0.401	ng/ul	99
13) Anthracene	16.95	178	4778	0.417	ng/ul	96
15) Fluoranthene	18.89	202	6202	0.387	ng/ul	97
17) Pyrene	19.25	202	6327	0.418	ng/ul	99
18) Benzo(a)anthracene	21.02	228	5333	0.402	ng/ul	99
19) Chrysene	21.07	228	5978	0.408	ng/ul	100
21) Benzo(b)fluoranthene	22.52	252	5827	0.414	ng/ul	96
22) Benzo(k)fluoranthene	22.56	252	6176	0.397	ng/ul	95
23) Benzo(a)pyrene	23.05	252	5349	0.409	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.20	276	6092	0.355	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.21	278	4875	0.359	ng/ul	96
26) Benzo(g,h,i)perylene	25.82	276	5245	0.338	ng/ul	99

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

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