

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012840.D
 Acq On : 03 Dec 2020 03:57
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Dec 03 04:30:13 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 : Wed Dec 02 12:19:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	718	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	2822	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1605	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	3261	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3043	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	3115	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	1737	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	3283	0.38	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.23	128	3588	0.422	ng/ul	99
5) 2-Methylnaphthalene	11.86	142	2466	0.455	ng/ul	99
7) Acenaphthylene	13.79	152	3527	0.451	ng/ul	99
8) Acenaphthene	14.14	153	2694	0.408	ng/ul	99
9) Fluorene	15.13	166	2981	0.411	ng/ul	99
11) Pentachlorophenol	16.49	266	473	0.457	ng/ul	96
12) Phenanthrene	16.87	178	4698	0.414	ng/ul	99
13) Anthracene	16.97	178	4228	0.429	ng/ul	97
15) Fluoranthene	18.90	202	5774	0.441	ng/ul	98
17) Pyrene	19.27	202	5923	0.435	ng/ul	100
18) Benzo(a)anthracene	21.03	228	5844	0.536	ng/ul	98
19) Chrysene	21.08	228	6809	0.500	ng/ul	99
21) Benzo(b)fluoranthene	22.53	252	6927	0.483	ng/ul	94
22) Benzo(k)fluoranthene	22.57	252	7693	0.466	ng/ul	94
23) Benzo(a)pyrene	23.06	252	6627	0.491	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.22	276	8768	0.497	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.23	278	6930	0.489	ng/ul	95
26) Benzo(g,h,i)perylene	25.85	276	7493	0.472	ng/ul	96

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

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