

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012199.D
 Acq On : 14 Oct 2020 17:55
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Oct 14 18:26:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 15:53:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1279	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5077	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2937	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5878	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	5093	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	5437	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2906	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	6202	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	5746	0.392	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	3671	0.394	ng/ul	99
7) Acenaphthylene	13.99	152	4845	0.380	ng/ul	99
8) Acenaphthene	14.33	153	4126	0.383	ng/ul	99
9) Fluorene	15.32	166	4586	0.383	ng/ul	100
11) Pentachlorophenol	16.67	266	659	0.388	ng/ul	99
12) Phenanthrene	17.06	178	7255	0.387	ng/ul	100
13) Anthracene	17.15	178	6224	0.387	ng/ul	99
15) Fluoranthene	19.08	202	8388	0.383	ng/ul	99
17) Pyrene	19.44	202	8802	0.383	ng/ul	99
18) Benzo(a)anthracene	21.20	228	7120	0.377	ng/ul	99
19) Chrysene	21.25	228	8859	0.385	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	8533	0.389	ng/ul	99
22) Benzo(k)fluoranthene	22.79	252	9477	0.390	ng/ul	98
23) Benzo(a)pyrene	23.31	252	8210	0.384	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.59	276	10730	0.391	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	8578	0.391	ng/ul	98
26) Benzo(g,h,i)perylene	26.25	276	9856	0.392	ng/ul	98

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

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