

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101420\
 Data File : BN012195.D
 Acq On : 14 Oct 2020 15:19
 Operator : CG/JU
 Sample : SSTD3.203
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.203

Quant Time: Oct 14 15:51:21 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 14:09:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1239	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5019	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2820	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5858	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	5313	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	5345	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	24979	3.17	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	52832	3.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	44490	2.995	ng/ul	95
5) 2-Methylnaphthalene	12.07	142	29831	3.012	ng/ul	99
7) Acenaphthylene	13.98	152	41257	3.034	ng/ul	97
8) Acenaphthene	14.32	153	32767	3.003	ng/ul	99
9) Fluorene	15.32	166	37811	2.998	ng/ul	98
11) Pentachlorophenol	16.67	266	5917	3.347	ng/ul	99
12) Phenanthrene	17.06	178	60008	3.046	ng/ul	97
13) Anthracene	17.14	178	55994	3.192	ng/ul	96
15) Fluoranthene	19.08	202	71383	3.260	ng/ul	96
17) Pyrene	19.44	202	72113	3.056	ng/ul	99
18) Benzo(a)anthracene	21.20	228	65962	3.193	ng/ul	98
19) Chrysene	21.25	228	71032	3.079	ng/ul	97
21) Benzo(b)fluoranthene	22.75	252	73428	2.931	ng/ul	84
22) Benzo(k)fluoranthene	22.79	252	76479	2.976	ng/ul#	86
23) Benzo(a)pyrene	23.31	252	67947	3.142	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.59	276	89618	3.153	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	72902	3.141	ng/ul#	87
26) Benzo(g,h,i)perylene	26.25	276	78279	3.152	ng/ul	95

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

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