

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN112020\
 Data File : BN012660.D
 Acq On : 20 Nov 2020 14:50
 Operator : CG/JU
 Sample : SSTD3.204
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.204

Quant Time: Nov 20 15:21:27 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 : Fri Nov 20 13:42:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	704	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	2650	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	1474	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3035	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3083	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	2731	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	12901	3.28	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	26532	3.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	24838	3.247	ng/ul#	94
5) 2-Methylnaphthalene	11.89	142	16634	3.424	ng/ul	99
7) Acenaphthylene	13.82	152	23333	3.649	ng/ul	96
8) Acenaphthene	14.16	153	18634	3.448	ng/ul	98
9) Fluorene	15.16	166	21435	3.563	ng/ul	99
11) Pentachlorophenol	16.51	266	3172	3.617	ng/ul	97
12) Phenanthrene	16.89	178	33883	3.502	ng/ul	96
13) Anthracene	16.99	178	31676	3.817	ng/ul	94
15) Fluoranthene	18.92	202	40763	3.603	ng/ul	96
17) Pyrene	19.29	202	40292	2.899	ng/ul	98
18) Benzo(a)anthracene	21.05	228	36498	3.196	ng/ul	96
19) Chrysene	21.10	228	39747	2.857	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	41060	3.722	ng/ul	87
22) Benzo(k)fluoranthene	22.60	252	44673	3.659	ng/ul#	88
23) Benzo(a)pyrene	23.09	252	38114	3.552	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.27	276	51293	3.721	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.28	278	41755	3.786	ng/ul#	88
26) Benzo(g,h,i)perylene	25.90	276	44232	3.501	ng/ul	94

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

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