

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009335.D
 Acq On : 10 Jan 2020 15:28
 Operator : JU
 Sample : SSTD1.601
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.601

Quant Time: Jan 10 16:00:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 14:47:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1477	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	5288	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2860	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	6406	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	5690	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	5334	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	13864	1.48	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	31307	1.66	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	25489	1.589	ng/ul	97
5) 2-Methylnaphthalene	11.90	142	17673	1.458	ng/ul	100
7) Acenaphthylene	13.82	152	24195	1.611	ng/ul	99
8) Acenaphthene	14.17	153	18946	1.607	ng/ul	100
9) Fluorene	15.17	166	22038	1.581	ng/ul	100
11) Pentachlorophenol	16.53	266	2966	1.737	ng/ul	98
12) Phenanthrene	16.90	178	35803	1.607	ng/ul	97
13) Anthracene	17.00	178	32749	1.639	ng/ul	97
15) Fluoranthene	18.93	202	42458	1.666	ng/ul	97
17) Pyrene	19.29	202	41901	1.443	ng/ul	99
18) Benzo(a)anthracene	21.05	228	39856	1.577	ng/ul	99
19) Chrysene	21.10	228	40731	1.604	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	39742	1.596	ng/ul	92
22) Benzo(k)fluoranthene	22.60	252	41779	1.664	ng/ul#	88
23) Benzo(a)pyrene	23.10	252	37080	1.699	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.26	276	44012	1.852	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.27	278	35180	1.859	ng/ul	94
26) Benzo(g,h,i)perylene	25.89	276	37702	1.883	ng/ul	96

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

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