

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011734.D
 Acq On : 08 Sep 2020 13:27
 Operator : CG/JU
 Sample : SSTD1.601
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.601

Quant Time: Sep 08 13:58:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 12:47:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1510	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	5569	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3141	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	6681	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5686	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	4975	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	14516	1.84	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	29263	1.56	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	26173	1.612	ng/ul	99
5) 2-Methylnaphthalene	12.19	142	17921	1.746	ng/ul	100
7) Acenaphthylene	14.09	152	24857	1.572	ng/ul	98
8) Acenaphthene	14.44	153	19655	1.597	ng/ul	99
9) Fluorene	15.43	166	23126	1.594	ng/ul	99
11) Pentachlorophenol	16.77	266	3252	2.050	ng/ul	99
12) Phenanthrene	17.16	178	36880	1.632	ng/ul	98
13) Anthracene	17.25	178	33998	1.829	ng/ul	99
15) Fluoranthene	19.18	202	42133	1.621	ng/ul	98
17) Pyrene	19.55	202	40963	1.843	ng/ul	99
18) Benzo(a)anthracene	21.30	228	36939	1.936	ng/ul	99
19) Chrysene	21.35	228	39659	1.541	ng/ul	98
21) Benzo(b)fluoranthene	22.89	252	39187	1.812	ng/ul	94
22) Benzo(k)fluoranthene	22.93	252	39582	1.656	ng/ul#	94
23) Benzo(a)pyrene	23.47	252	33304	1.833	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.83	276	44042	1.793	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	36469	1.821	ng/ul	96
26) Benzo(g,h,i)perylene	26.52	276	38020	1.653	ng/ul	100

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

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