

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011731.D
 Acq On : 08 Sep 2020 11:38
 Operator : CG/JU
 Sample : SST00.204
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.204

Quant Time: Sep 08 12:55:22 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	1762	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	6296	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3336	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7180	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5862	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	5402	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	1864	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	3514	0.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	3681	0.201	ng/ul	98
5) 2-Methylnaphthalene	12.19	142	2409	0.208	ng/ul	98
7) Acenaphthylene	14.09	152	3108	0.185	ng/ul	98
8) Acenaphthene	14.44	153	2521	0.193	ng/ul	98
9) Fluorene	15.43	166	2884	0.187	ng/ul	99
11) Pentachlorophenol	16.77	266	403	0.236	ng/ul	97
12) Phenanthrene	17.16	178	4629	0.191	ng/ul	99
13) Anthracene	17.25	178	4129	0.207	ng/ul	97
15) Fluoranthene	19.18	202	5047	0.181	ng/ul	97
17) Pyrene	19.55	202	5062	0.221	ng/ul	98
18) Benzo(a)anthracene	21.30	228	4309	0.219	ng/ul	99
19) Chrysene	21.35	228	4940	0.186	ng/ul	100
21) Benzo(b)fluoranthene	22.89	252	4871	0.207	ng/ul	88
22) Benzo(k)fluoranthene	22.94	252	4970	0.191	ng/ul#	91
23) Benzo(a)pyrene	23.47	252	4082	0.207	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.83	276	5391	0.202	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	4396	0.202	ng/ul#	92
26) Benzo(g,h,i)perylene	26.52	276	4805	0.192	ng/ul	94

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

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