

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN090820\
 Data File : BN011735.D
 Acq On : 08 Sep 2020 14:03
 Operator : CG/JU
 Sample : SSTD3.202
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.202

Quant Time: Sep 08 14:32:48 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	1624	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	5929	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3222	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	6680	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5914	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	5196	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	29749	3.55	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	58575	3.12	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	54203	3.136	ng/ul	97
5) 2-Methylnaphthalene	12.19	142	36667	3.356	ng/ul	100
7) Acenaphthylene	14.09	152	50754	3.129	ng/ul	99
8) Acenaphthene	14.44	153	39499	3.128	ng/ul	98
9) Fluorene	15.43	166	46316	3.112	ng/ul	98
11) Pentachlorophenol	16.77	266	6674	4.208	ng/ul	99
12) Phenanthrene	17.16	178	72279	3.198	ng/ul	98
13) Anthracene	17.25	178	68811	3.702	ng/ul	99
15) Fluoranthene	19.18	202	83854	3.226	ng/ul	99
17) Pyrene	19.55	202	82244	3.558	ng/ul	98
18) Benzo(a)anthracene	21.30	228	76259	3.843	ng/ul	100
19) Chrysene	21.35	228	80003	2.990	ng/ul	99
21) Benzo(b)fluoranthene	22.89	252	80281	3.555	ng/ul	94
22) Benzo(k)fluoranthene	22.93	252	81948	3.282	ng/ul#	92
23) Benzo(a)pyrene	23.47	252	69899	3.683	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.83	276	94079	3.666	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	77881	3.723	ng/ul	95
26) Benzo(g,h,i)perylene	26.52	276	80544	3.353	ng/ul	98

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

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