

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072020\
 Data File : BN011202.D
 Acq On : 20 Jul 2020 09:58
 Operator : CG/JU
 Sample : SST00.206
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.206

Quant Time: Jul 20 11:28:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 20 11:25:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	1067	0.40	ng/ul	0.00
2) Naphthalene-d8	10.25	136	3767	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	2256	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	4782	0.40	ng/ul	0.00
16) Chrysene-d12	21.10	240	3144	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	2846	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	1747	0.20	ng/ul	0.01
14) Fluoranthene-d10	18.92	212	3956	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.30	128	2198	0.198	ng/ul#	93
5) 2-Methylnaphthalene	11.92	142	1392	0.200	ng/ul	98
7) Acenaphthylene	13.84	152	2131	0.212	ng/ul	95
8) Acenaphthene	14.19	153	1680	0.199	ng/ul	99
9) Fluorene	15.19	166	1883	0.203	ng/ul	97
11) Pentachlorophenol	16.56	266	177	0.158	ng/ul	93
12) Phenanthrene	16.92	178	3066	0.197	ng/ul	97
13) Anthracene	17.02	178	2398	0.187	ng/ul	94
15) Fluoranthene	18.95	202	3347	0.192	ng/ul	97
17) Pyrene	19.32	202	3405	0.249	ng/ul#	96
18) Benzo(a)anthracene	21.09	228	2063	0.185	ng/ul	96
19) Chrysene	21.13	228	2972	0.209	ng/ul	97
21) Benzo(b)fluoranthene	22.61	252	2320	0.183	ng/ul	85
22) Benzo(k)fluoranthene	22.65	252	2783	0.207	ng/ul#	83
23) Benzo(a)pyrene	23.15	252	2017	0.188	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.33	276	2260	0.154	ng/ul	98
25) Dibenzo(a,h)anthracene	25.35	278	1680	0.140	ng/ul#	80
26) Benzo(g,h,i)perylene	25.95	276	2372	0.175	ng/ul	93

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

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