

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073020\
 Data File : BN011311.D
 Acq On : 30 Jul 2020 22:53
 Operator : CG/JU
 Sample : L3388-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD6

Quant Time: Jul 31 03:01:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 02:58:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	858	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	3550	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2301	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5449	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	5742	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	6581	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	1210	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	3699	0.14	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.52	266	82	0.061	ng/ul#	87
12) Phenanthrene	16.89	178	615	0.034	ng/ul#	82
13) Anthracene	16.99	178	379	0.024	ng/ul#	61
15) Fluoranthene	18.93	202	1754	0.079	ng/ul	88
17) Pyrene	19.29	202	1800	0.069	ng/ul#	91
18) Benzo(a)anthracene	21.06	228	1427	0.065	ng/ul#	91
19) Chrysene	21.11	228	1476	0.058	ng/ul#	93
21) Benzo(b)fluoranthene	22.57	252	1707	0.062	ng/ul#	45
22) Benzo(k)fluoranthene	22.61	252	1614	0.052	ng/ul#	28
23) Benzo(a)pyrene	23.11	252	1313	0.055	ng/ul#	42
24) Indeno(1,2,3-cd)pyrene	25.27	276	1864	0.074	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.28	278	1377	0.069	ng/ul#	66
26) Benzo(g,h,i)perylene	25.90	276	1761	0.078	ng/ul#	78

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

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