

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012804.D
 Acq On : 02 Dec 2020 05:05
 Operator : CG/JU
 Sample : L4793-21
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B54

Manual Integrations
 APPROVED

mohammad
 12/2/2020 2:28:35 PM

Quant Time: Dec 02 07:07:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 05:08:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	945	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	3903	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	2269	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	4828	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	4003	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	4369	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	672	0.12	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	1627	0.13	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.87	178	1517	0.090	ng/ul	95
15) Fluoranthene	18.90	202	5003	0.258	ng/ul	98
17) Pyrene	19.27	202	5062	0.283	ng/ul	97
18) Benzo(a)anthracene	21.04	228	2175	0.152	ng/ul	96
19) Chrysene	21.09	228	3005	0.168	ng/ul	95
21) Benzo(b)fluoranthene	22.54	252	4494m	0.224	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	1629m	0.070	ng/ul	
23) Benzo(a)pyrene	23.07	252	2729	0.144	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.23	276	2230	0.090	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.23	278	498	0.025	ng/ul#	22
26) Benzo(g,h,i)perylene	25.86	276	2131	0.096	ng/ul#	91

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

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