

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012212.D
 Acq On : 15 Oct 2020 01:44
 Operator : CG/JU
 Sample : L4166-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 15 02:27:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 00:46:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1261	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5077	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3014	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	6085m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	5035	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	5943	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1869	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	3967	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.98	152	321	0.025	ng/ul#	62
13) Anthracene	17.14	178	337	0.020	ng/ul#	64
15) Fluoranthene	19.07	202	715	0.032	ng/ul	80
17) Pyrene	19.44	202	866	0.038	ng/ul#	70
18) Benzo(a)anthracene	21.20	228	732	0.039	ng/ul#	78
19) Chrysene	21.24	228	1472	0.065	ng/ul	95
21) Benzo(b)fluoranthene	22.74	252	2201m	0.092	ng/ul	
22) Benzo(k)fluoranthene	22.78	252	909m	0.034	ng/ul	
23) Benzo(a)pyrene	23.30	252	1458	0.062	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	25.57	276	985	0.033	ng/ul#	100
26) Benzo(g,h,i)perylene	26.24	276	991	0.036	ng/ul#	68

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

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