

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102219\
 Data File : BN008365.D
 Acq On : 24 Oct 2019 00:30
 Operator : JU
 Sample : K5235-03
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPB1

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:07:14 AM

Quant Time: Oct 24 02:22:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1862	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	9077	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5863	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12974	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	11262	0.40	ng/ul	-0.01
20) Perylene-d12	23.33	264	9980	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2124	0.14	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	6394	0.14	ng/ul	-0.01
Target Compounds				Qvalue		
12) Phenanthrene	17.00	178	3008	0.079	ng/ul	96
15) Fluoranthene	19.03	202	7615	0.152	ng/ul	94
17) Pyrene	19.39	202	7285m	0.162	ng/ul	
18) Benzo(a)anthracene	21.15	228	4487	0.110	ng/ul	93
19) Chrysene	21.20	228	4468	0.089	ng/ul	94
21) Benzo(b)fluoranthene	22.70	252	5301m	0.163	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	1626m	0.044	ng/ul	
23) Benzo(a)pyrene	23.24	252	3626	0.106	ng/ul#	56
24) Indeno(1,2,3-cd)pyrene	25.49	276	2041	0.051	ng/ul#	89
26) Benzo(g,h,i)perylene	26.14	276	2048	0.055	ng/ul#	81

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

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