

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008197.D
 Acq On : 16 Oct 2019 14:17
 Operator : JU
 Sample : K5175-20
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP92

Manual Integrations
 APPROVED

mohammad
 10/18/2019 9:13:23 AM

Quant Time: Oct 16 18:11:11 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 16 17:43:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2530	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	12375	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7999	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	17652m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	16268	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	13908	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	2309	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	7101	0.12	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.03	178	9651	0.187	ng/ul	100
13) Anthracene	17.12	178	1955	0.043	ng/ul#	91
15) Fluoranthene	19.06	202	23440	0.343	ng/ul	98
17) Pyrene	19.42	202	23820	0.368	ng/ul	98
18) Benzo(a)anthracene	21.17	228	13543	0.229	ng/ul	95
19) Chrysene	21.23	228	13483	0.186	ng/ul	97
21) Benzo(b)fluoranthene	22.73	252	14427m	0.319	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	5752m	0.112	ng/ul	
23) Benzo(a)pyrene	23.28	252	10583	0.222	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.54	276	5203	0.093	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.54	278	1186	0.027	ng/ul#	26
26) Benzo(g,h,i)perylene	26.20	276	5357	0.104	ng/ul#	86

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

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