

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008194.D
 Acq On : 16 Oct 2019 12:29
 Operator : JU
 Sample : K5175-17
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPG4

Manual Integrations
 APPROVED

mohammad
 10/18/2019 9:13:15 AM

Quant Time: Oct 16 18:02:25 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	2976	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	14447	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	9137	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	20292m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	18767	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	15826	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	6160	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	18669	0.27	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.03	178	7286	0.123	ng/ul	99
13) Anthracene	17.12	178	1499	0.029	ng/ul#	86
15) Fluoranthene	19.06	202	19526	0.249	ng/ul	97
17) Pyrene	19.42	202	18907	0.253	ng/ul#	94
18) Benzo(a)anthracene	21.18	228	9782	0.143	ng/ul	94
19) Chrysene	21.23	228	11581	0.138	ng/ul	95
21) Benzo(b)fluoranthene	22.73	252	12777m	0.248	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	4650m	0.080	ng/ul	
23) Benzo(a)pyrene	23.28	252	8142	0.150	ng/ul#	55
24) Indeno(1,2,3-cd)pyrene	25.54	276	4382	0.069	ng/ul#	91
26) Benzo(g,h,i)perylene	26.20	276	3792	0.065	ng/ul#	77

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

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