

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008188.D
 Acq On : 16 Oct 2019 00:01
 Operator : JU
 Sample : K5175-14
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPG1

Manual Integrations
 APPROVED

mohammad
 10/16/2019 4:09:02 PM

Quant Time: Oct 16 05:07:40 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 : Tue Oct 15 15:13:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2245	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.38	136	10796	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.25	164	6899	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.00	188	15934m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	16879m	0.40	ng/ul	-0.01
20) Perylene-d12	23.37	264	12508m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5143	0.28	ng/ul	-0.01
14) Fluoranthene-d10	19.03	212	16847m	0.31	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	816	0.027	ng/ul#	76
5) 2-Methylnaphthalene	12.06	142	472	0.022	ng/ul	98
7) Acenaphthylene	13.96	152	2317	0.071	ng/ul#	92
8) Acenaphthene	14.31	153	630	0.025	ng/ul	96
9) Fluorene	15.30	166	678	0.023	ng/ul#	91
12) Phenanthrene	17.04	178	13849	0.297	ng/ul	100
13) Anthracene	17.13	178	2831m	0.069	ng/ul	
15) Fluoranthene	19.06	202	36194m	0.587	ng/ul	
17) Pyrene	19.42	202	32787m	0.488	ng/ul	
18) Benzo(a)anthracene	21.18	228	21757m	0.355	ng/ul	
19) Chrysene	21.23	228	21558	0.286	ng/ul	98
21) Benzo(b)fluoranthene	22.73	252	26008m	0.639	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	8588m	0.186	ng/ul	
23) Benzo(a)pyrene	23.28	252	17291	0.404	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.54	276	10049	0.199	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.55	278	2275	0.057	ng/ul#	62
26) Benzo(a,h,i)perylene	26.19	276	7175	0.155	ng/ul#	92

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

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