

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008184.D
 Acq On : 15 Oct 2019 21:37
 Operator : JU
 Sample : K5175-12
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPF9

Manual Integrations
 APPROVED

mohammad
 10/16/2019 4:08:44 PM

Quant Time: Oct 16 04:47:35 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	2215	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.38	136	10781	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.25	164	6875	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	15671m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	16289m	0.40	ng/ul	-0.01
20) Perylene-d12	23.38	264	13249m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5863	0.32	ng/ul	-0.01
14) Fluoranthene-d10	19.03	212	19114m	0.35	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	1619	0.053	ng/ul#	87
5) 2-Methylnaphthalene	12.06	142	823	0.039	ng/ul	97
7) Acenaphthylene	13.96	152	7809	0.238	ng/ul	98
8) Acenaphthene	14.31	153	3542	0.139	ng/ul	98
9) Fluorene	15.30	166	4301	0.149	ng/ul#	96
12) Phenanthrene	17.04	178	101344m	2.209	ng/ul	
13) Anthracene	17.13	178	21812m	0.540	ng/ul	
15) Fluoranthene	19.06	202	269725m	4.451	ng/ul	
17) Pyrene	19.42	202	289757	4.468	ng/ul	98
18) Benzo(a)anthracene	21.18	228	176984m	2.989	ng/ul	
19) Chrysene	21.23	228	183950	2.531	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	189240m	4.388	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	61350m	1.256	ng/ul	
23) Benzo(a)pyrene	23.28	252	136524m	3.008	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.54	276	69738	1.305	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.55	278	18814	0.445	ng/ul	95
26) Benzo(a,h,i)perylene	26.20	276	57991	1.183	ng/ul	99

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

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