

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008182.D
 Acq On : 15 Oct 2019 20:26
 Operator : JU
 Sample : K5175-10
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPF7

Manual Integrations
 APPROVED

mohammad
 10/16/2019 4:08:40 PM

Quant Time: Oct 16 04:22:50 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	2449	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.38	136	11912	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.25	164	7414	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	16983m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	17179m	0.40	ng/ul	-0.01
20) Perylene-d12	23.38	264	15304m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4171	0.21	ng/ul	-0.01
14) Fluoranthene-d10	19.03	212	14019m	0.24	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.43	128	5346	0.158	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	4179	0.180	ng/ul	99
7) Acenaphthylene	13.96	152	13933	0.395	ng/ul	99
8) Acenaphthene	14.31	153	8078	0.294	ng/ul	99
9) Fluorene	15.30	166	8499	0.273	ng/ul#	95
12) Phenanthrene	17.03	178	197869m	3.979	ng/ul	
13) Anthracene	17.13	178	38657m	0.883	ng/ul	
15) Fluoranthene	19.06	202	398088m	6.062	ng/ul	
17) Pyrene	19.43	202	469119	6.858	ng/ul	99
18) Benzo(a)anthracene	21.18	228	304022m	4.868	ng/ul	
19) Chrysene	21.23	228	321089	4.189	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	294110m	5.903	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	102829m	1.822	ng/ul	
23) Benzo(a)pyrene	23.28	252	236438	4.510	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.54	276	111656	1.808	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	33131	0.679	ng/ul	96
26) Benzo(a,h,i)perylene	26.20	276	100414	1.773	ng/ul	99

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

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