

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006504.D
 Acq On : 22 Jun 2019 23:51
 Operator : HP/JU
 Sample : K3360-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Z0

Manual Integrations
 APPROVED

mohammad
 6/25/2019 8:44:56 AM

Quant Time: Jun 24 09:21:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5007	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	19849	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	10957	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	22483m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.25	240	14021m	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	10949m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	5565	0.18	ng/ul	-0.01
14) Fluoranthene-d10	19.08	212	13086m	0.22	ng/ul	-0.01
Target Compounds						
3) Naphthalene	10.46	128	2345	0.043	ng/ul#	89
5) 2-Methylnaphthalene	12.08	142	771	0.020	ng/ul	92
12) Phenanthrene	17.08	178	3246	0.045	ng/ul	93
15) Fluoranthene	19.11	202	8487m	0.106	ng/ul	
17) Pyrene	19.47	202	8903	0.121	ng/ul	96
18) Benzo(a)anthracene	21.23	228	4373	0.071	ng/ul	96
19) Chrysene	21.29	228	5206	0.089	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	5620	0.123	ng/ul	83
22) Benzo(k)fluoranthene	22.86	252	3871m	0.087	ng/ul	
23) Benzo(a)pyrene	23.38	252	3722m	0.088	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.71	276	3733m	0.081	ng/ul	
25) Dibenzo(a,h)anthracene	25.72	278	2220m	0.060	ng/ul	
26) Benzo(g,h,i)perylene	26.39	276	3063m	0.079	ng/ul	

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

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