

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006556.D
 Acq On : 25 Jun 2019 05:51
 Operator : HP/JU
 Sample : K3388-03
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4216

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:21:24 PM

Quant Time: Jun 25 07:45:41 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 : Tue Jun 25 01:51:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5254	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	19638	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9740	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	18320m	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	9854m	0.40	ng/ul	-0.01
20) Perylene-d12	23.49	264	9796m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	5947	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	11538m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	1243	0.023	ng/ul#	79
7) Acenaphthylene	14.00	152	1618	0.030	ng/ul#	78
12) Phenanthrene	17.08	178	12995	0.221	ng/ul	97
13) Anthracene	17.17	178	2218m	0.039	ng/ul	
15) Fluoranthene	19.11	202	31079m	0.479	ng/ul	
17) Pyrene	19.48	202	25628	0.494	ng/ul	99
18) Benzo(a)anthracene	21.24	228	9571m	0.220	ng/ul	
19) Chrysene	21.29	228	11364m	0.278	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	16448m	0.403	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	5517m	0.139	ng/ul	
23) Benzo(a)pyrene	23.39	252	10030	0.264	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.72	276	8027	0.195	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.72	278	1944	0.059	ng/ul#	1
26) Benzo(g,h,i)perylene	26.41	276	6809	0.197	ng/ul#	75

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

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