

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006514.D
 Acq On : 23 Jun 2019 07:19
 Operator : HP/JU
 Sample : K3360-12
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Z2

Manual Integrations
 APPROVED

mohammad
 6/25/2019 8:45:27 AM

Quant Time: Jun 24 09:09:17 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	5304	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	20329	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	10747	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.04	188	21638m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.25	240	12676m	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	11062m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	6538	0.21	ng/ul	-0.02
14) Fluoranthene-d10	19.08	212	13272m	0.23	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	2141	0.038	ng/ul#	87
7) Acenaphthylene	14.00	152	7421	0.124	ng/ul	97
12) Phenanthrene	17.08	178	24354m	0.351	ng/ul	
13) Anthracene	17.17	178	6724m	0.101	ng/ul	
15) Fluoranthene	19.10	202	120207m	1.567	ng/ul	
17) Pyrene	19.47	202	102489	1.535	ng/ul	99
18) Benzo(a)anthracene	21.24	228	44720	0.798	ng/ul	99
19) Chrysene	21.29	228	50266	0.955	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	59577m	1.292	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	23936m	0.533	ng/ul	
23) Benzo(a)pyrene	23.38	252	36827m	0.857	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.71	276	27422m	0.590	ng/ul	
25) Dibenzo(a,h)anthracene	25.72	278	6599m	0.178	ng/ul	
26) Benzo(g,h,i)perylene	26.40	276	22249m	0.571	ng/ul	

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

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