

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006496.D
 Acq On : 22 Jun 2019 19:15
 Operator : HP/JU
 Sample : K3335-07RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4234RE

Manual Integrations
 APPROVED

mohammad
 6/25/2019 8:44:31 AM

Quant Time: Jun 24 08:36:37 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	4533	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	18134	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	9956	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	294686	0.40	ng/ul	0.01
16) Chrysene-d12	21.26	240	14654	0.40	ng/ul	-0.01
20) Perylene-d12	23.48	264	11022m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	7069	0.25	ng/ul	-0.02
14) Fluoranthene-d10	19.08	212	17413	0.02	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.46	128	3629	0.073	ng/ul#	93
5) 2-Methylnaphthalene	12.08	142	1269	0.036	ng/ul	99
7) Acenaphthylene	14.00	152	7605	0.137	ng/ul	98
8) Acenaphthene	14.34	153	3132	0.082	ng/ul	98
9) Fluorene	15.34	166	4401	0.098	ng/ul#	97
12) Phenanthrene	17.08	178	84856	0.090	ng/ul	99
13) Anthracene	17.17	178	20426	0.023	ng/ul	99
15) Fluoranthene	19.11	202	178260	0.171	ng/ul	99
17) Pyrene	19.48	202	163858	2.123	ng/ul	99
18) Benzo(a)anthracene	21.24	228	70068	1.081	ng/ul	100
19) Chrysene	21.29	228	68019	1.118	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	77990m	1.697	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	28265m	0.631	ng/ul	
23) Benzo(a)pyrene	23.39	252	51881m	1.212	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.71	276	31864m	0.688	ng/ul	
25) Dibenzo(a,h)anthracene	25.72	278	8393m	0.227	ng/ul	
26) Benzo(g,h,i)perylene	26.40	276	25036m	0.645	ng/ul	

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

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