

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009050.D
 Acq On : 20 Dec 2019 03:48
 Operator : JU
 Sample : K6171-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BS6

Manual Integrations
 APPROVED

Sohil
 12/20/2019 4:30:00 PM

Quant Time: Dec 20 06:24:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 19 15:54:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	3375	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	13286	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.17	164	7727	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	16070	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	9439	0.40	ng/ul	0.00
20) Perylene-d12	23.25	264	8107	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	4460	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.95	212	10202	0.22	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.34	128	4609	0.114	ng/ul	97
5) 2-Methylnaphthalene	11.96	142	4002	0.131	ng/ul	99
7) Acenaphthylene	13.88	152	14457	0.356	ng/ul	99
8) Acenaphthene	14.23	153	1819	0.057	ng/ul	97
9) Fluorene	15.22	166	4985	0.132	ng/ul#	98
12) Phenanthrene	16.95	178	138304	2.475	ng/ul	100
13) Anthracene	17.05	178	30284	0.604	ng/ul	99
15) Fluoranthene	18.98	202	545980	8.543	ng/ul	98
17) Pyrene	19.34	202	405888	8.429	ng/ul	99
18) Benzo(a)anthracene	21.10	228	181005	4.318	ng/ul	97
19) Chrysene	21.15	228	158948	3.773	ng/ul	97
21) Benzo(b)fluoranthene	22.62	252	187576m	4.955	ng/ul	
22) Benzo(k)fluoranthene	22.66	252	65497m	1.716	ng/ul	
23) Benzo(a)pyrene	23.16	252	115329	3.477	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.35	276	78093	2.162	ng/ul	91
25) Dibenzo(a,h)anthracene	25.36	278	17895	0.622	ng/ul	92
26) Benzo(a,h,i)perylene	25.99	276	73348	2.410	ng/ul	99

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

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