

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009052.D
 Acq On : 20 Dec 2019 04:59
 Operator : JU
 Sample : K6171-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BS9

Manual Integrations
 APPROVED

Sohil
 12/20/2019 4:30:01 PM

Quant Time: Dec 20 06:29:49 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	3623	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	13971	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.17	164	7991	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	15964m	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	8742	0.40	ng/ul	0.00
20) Perylene-d12	23.25	264	8120	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	5949	0.24	ng/ul	0.00
14) Fluoranthene-d10	18.94	212	11360	0.24	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.34	128	6416	0.151	ng/ul	98
5) 2-Methylnaphthalene	11.96	142	5009	0.156	ng/ul	99
7) Acenaphthylene	13.88	152	5713	0.136	ng/ul	96
8) Acenaphthene	14.23	153	2558	0.078	ng/ul	99
9) Fluorene	15.22	166	6348	0.163	ng/ul#	98
12) Phenanthrene	16.95	178	134913	2.430	ng/ul	100
13) Anthracene	17.05	178	20962	0.421	ng/ul	99
15) Fluoranthene	18.98	202	286901	4.519	ng/ul	98
17) Pyrene	19.34	202	214551	4.811	ng/ul	96
18) Benzo(a)anthracene	21.10	228	84301	2.172	ng/ul	96
19) Chrysene	21.15	228	95044	2.436	ng/ul	96
21) Benzo(b)fluoranthene	22.62	252	115402m	3.044	ng/ul	
22) Benzo(k)fluoranthene	22.66	252	35306m	0.923	ng/ul	
23) Benzo(a)pyrene	23.16	252	74316	2.237	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.35	276	54234	1.499	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.36	278	12978	0.451	ng/ul#	79
26) Benzo(a,h,i)perylene	25.99	276	58261	1.911	ng/ul	99

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

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