

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009056.D
 Acq On : 20 Dec 2019 07:21
 Operator : JU
 Sample : K6171-12
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BW1

Manual Integrations
 APPROVED

Sohil
 12/20/2019 4:30:06 PM

Quant Time: Dec 20 11:01:39 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	4261	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	16534	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.17	164	8996	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	17482	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	9000	0.40	ng/ul	0.00
20) Perylene-d12	23.26	264	8633	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	5135	0.17	ng/ul	0.00
14) Fluoranthene-d10	18.95	212	9026	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.34	128	4662	0.093	ng/ul	96
5) 2-Methylnaphthalene	11.96	142	4579	0.121	ng/ul	100
7) Acenaphthylene	13.88	152	3297	0.070	ng/ul#	92
8) Acenaphthene	14.23	153	1419	0.038	ng/ul	97
9) Fluorene	15.22	166	3066	0.070	ng/ul#	95
12) Phenanthrene	16.95	178	68976	1.135	ng/ul	100
13) Anthracene	17.05	178	10322	0.189	ng/ul	98
15) Fluoranthene	18.98	202	140861	2.026	ng/ul	97
17) Pyrene	19.34	202	114745	2.499	ng/ul	99
18) Benzo(a)anthracene	21.10	228	52370	1.310	ng/ul#	87
19) Chrysene	21.15	228	63306	1.576	ng/ul#	92
21) Benzo(b)fluoranthene	22.63	252	74738m	1.854	ng/ul	
22) Benzo(k)fluoranthene	22.66	252	26446m	0.651	ng/ul	
23) Benzo(a)pyrene	23.16	252	58335	1.651	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.36	276	41875	1.088	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.36	278	10725	0.350	ng/ul#	61
26) Benzo(a,h,i)perylene	25.99	276	64434	1.988	ng/ul	99

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

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