

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008659.D
 Acq On : 20 Nov 2019 14:19
 Operator : JU
 Sample : K5784-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLNBI

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:05:57 PM

Quant Time: Nov 20 15:12:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 20 04:45:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	4082	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	15254	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8457	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	17278	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	14005	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	14251	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4172	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	9514	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	2361	0.051	ng/ul#	90
5) 2-Methylnaphthalene	12.04	142	1491	0.046	ng/ul	99
7) Acenaphthylene	13.96	152	2018	0.043	ng/ul#	86
8) Acenaphthene	14.30	153	2871	0.083	ng/ul	99
9) Fluorene	15.29	166	2329	0.057	ng/ul#	93
12) Phenanthrene	17.03	178	48633	0.835	ng/ul	98
13) Anthracene	17.11	178	9413	0.171	ng/ul	98
15) Fluoranthene	19.05	202	70195	0.981	ng/ul	96
17) Pyrene	19.41	202	78468	1.397	ng/ul	99
18) Benzo(a)anthracene	21.16	228	28710	0.510	ng/ul#	91
19) Chrysene	21.21	228	32447	0.587	ng/ul#	95
21) Benzo(b)fluoranthene	22.71	252	31847m	0.542	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	12183m	0.209	ng/ul	
23) Benzo(a)pyrene	23.25	252	27382	0.496	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.49	276	18015	0.279	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.49	278	4230	0.081	ng/ul#	9
26) Benzo(a,h,i)perylene	26.14	276	19718	0.365	ng/ul#	89

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

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