

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008660.D
 Acq On : 20 Nov 2019 14:55
 Operator : JU
 Sample : K5784-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLN2

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:06:00 PM

Quant Time: Nov 20 17:42:05 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	3736	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	14097	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7705	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	16143	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13609	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13954	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3354	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	8894	0.18	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.03	178	4772	0.088	ng/ul	98
15) Fluoranthene	19.05	202	10175	0.152	ng/ul	93
17) Pyrene	19.41	202	12629	0.231	ng/ul	97
18) Benzo(a)anthracene	21.16	228	5916	0.108	ng/ul#	90
19) Chrysene	21.21	228	6419	0.119	ng/ul#	89
21) Benzo(b)fluoranthene	22.71	252	8069m	0.140	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	3044m	0.053	ng/ul	
23) Benzo(a)pyrene	23.25	252	6193	0.115	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	25.48	276	4162	0.066	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.50	278	1054	0.021	ng/ul#	1
26) Benzo(g,h,i)perylene	26.14	276	4508	0.085	ng/ul#	70

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

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