

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112420\
 Data File : BN012715.D
 Acq On : 24 Nov 2020 14:16
 Operator : CG/JU
 Sample : L4751-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BG0

Manual Integrations
 APPROVED

mohammad
 11/25/2020 1:20:47 PM

Quant Time: Nov 24 14:51:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 24 11:59:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	882	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	3718	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2213	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	4428	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3603	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3841	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	731	0.14	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	1736	0.15	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.81	152	241	0.022	ng/ul#	44
12) Phenanthrene	16.89	178	1579	0.102	ng/ul	94
15) Fluoranthene	18.92	202	5274	0.296	ng/ul	98
17) Pyrene	19.28	202	4985	0.309	ng/ul#	91
18) Benzo(a)anthracene	21.05	228	2356	0.182	ng/ul	92
19) Chrysene	21.10	228	3010m	0.187	ng/ul	
21) Benzo(b)fluoranthene	22.56	252	4521m	0.256	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	1973m	0.097	ng/ul	
23) Benzo(a)pyrene	23.08	252	2803	0.169	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.26	276	2149	0.099	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.26	278	501	0.029	ng/ul#	3
26) Benzo(g,h,i)perylene	25.89	276	2030	0.104	ng/ul#	82

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

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