

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011385.D
 Acq On : 06 Aug 2020 18:38
 Operator : CG/JU
 Sample : L3419-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AF0

Manual Integrations
 APPROVED

mohammad
 8/7/2020 1:53:31 PM

Quant Time: Aug 07 02:09:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 07 02:04:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	986	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	3467	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1878	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	4168	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	4241	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4643	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	1134	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	2719	0.14	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.80	152	180	0.020	ng/ul#	47
9) Fluorene	15.15	166	208	0.025	ng/ul#	91
12) Phenanthrene	16.88	178	3795	0.275	ng/ul	98
13) Anthracene	16.97	178	280	0.023	ng/ul#	43
15) Fluoranthene	18.91	202	9494	0.558	ng/ul	99
17) Pyrene	19.27	202	7289	0.378	ng/ul	100
18) Benzo(a)anthracene	21.05	228	1809	0.111	ng/ul#	93
19) Chrysene	21.10	228	2000	0.107	ng/ul#	92
21) Benzo(b)fluoranthene	22.55	252	2263m	0.117	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	879m	0.040	ng/ul	
23) Benzo(a)pyrene	23.08	252	660	0.039	ng/ul#	6
24) Indeno(1,2,3-cd)pyrene	25.25	276	579	0.033	ng/ul#	82
26) Benzo(g,h,i)perylene	25.87	276	520	0.032	ng/ul#	55

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

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