

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013004.D
 Acq On : 09 Dec 2020 09:38
 Operator : CG/JU
 Sample : L4942-01DL 10X
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B60DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:26:08 PM

Quant Time: Dec 09 10:26:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1265	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.14	136	5106	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.05	164	2770	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5602	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4088	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3675	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	114	0.02	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	234	0.02	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	463	0.035	ng/ul#	68
8) Acenaphthene	14.11	153	509	0.047	ng/ul	97
9) Fluorene	15.11	166	397	0.032	ng/ul#	95
12) Phenanthrene	16.85	178	4928	0.262	ng/ul	99
13) Anthracene	16.94	178	857	0.054	ng/ul#	87
15) Fluoranthene	18.88	202	15637	0.701	ng/ul	99
17) Pyrene	19.25	202	13871	0.718	ng/ul	98
18) Benzo(a)anthracene	21.01	228	6484	0.383	ng/ul	99
19) Chrysene	21.07	228	7291	0.390	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	8847m	0.520	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	3339m	0.178	ng/ul	
23) Benzo(a)pyrene	23.04	252	5528	0.350	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.19	276	3616	0.174	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.19	278	827	0.050	ng/ul#	27
26) Benzo(g,h,i)perylene	25.82	276	3169	0.169	ng/ul#	87

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

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