

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012924.D
 Acq On : 06 Dec 2020 19:05
 Operator : CG/JU
 Sample : L4864-19
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B00

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:11:54 PM

Quant Time: Dec 07 03:51:20 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 : Mon Dec 07 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	2044	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	8531	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	4722	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	9968	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	7697	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	6837	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1573	0.12	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	3520	0.13	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	794	0.032	ng/ul#	93
5) 2-Methylnaphthalene	11.84	142	358	0.022	ng/ul	95
7) Acenaphthylene	13.77	152	589	0.026	ng/ul#	82
8) Acenaphthene	14.12	153	831	0.045	ng/ul	96
9) Fluorene	15.12	166	942	0.045	ng/ul#	97
12) Phenanthrene	16.85	178	15282	0.457	ng/ul	97
13) Anthracene	16.95	178	3136	0.110	ng/ul	98
15) Fluoranthene	18.89	202	21239	0.535	ng/ul	98
17) Pyrene	19.25	202	18508	0.509	ng/ul	96
18) Benzo(a)anthracene	21.02	228	8901	0.280	ng/ul	99
19) Chrysene	21.07	228	8644	0.246	ng/ul	97
21) Benzo(b)fluoranthene	22.52	252	10867m	0.344	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	4077m	0.117	ng/ul	
23) Benzo(a)pyrene	23.05	252	7100	0.242	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.19	276	4232	0.110	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.20	278	1102	0.036	ng/ul#	38
26) Benzo(a,h,i)perylene	25.82	276	4038	0.116	ng/ul#	91

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

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