

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012929.D
 Acq On : 06 Dec 2020 22:06
 Operator : CG/JU
 Sample : L4864-21
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B64

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:12:08 PM

Quant Time: Dec 07 04:08:52 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	1248	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	5176	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	2928	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	6681	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	5722	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	5476	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	987	0.13	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	2361	0.13	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	313	0.021	ng/ul#	67
7) Acenaphthylene	13.77	152	500	0.036	ng/ul#	79
12) Phenanthrene	16.85	178	12356	0.552	ng/ul	98
13) Anthracene	16.95	178	1760	0.092	ng/ul	96
15) Fluoranthene	18.88	202	59544	2.237	ng/ul	100
17) Pyrene	19.25	202	54064	2.000	ng/ul	98
18) Benzo(a)anthracene	21.02	228	23668	1.000	ng/ul	99
19) Chrysene	21.07	228	31090	1.189	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	47078m	1.858	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	13804m	0.493	ng/ul	
23) Benzo(a)pyrene	23.05	252	26882m	1.144	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	20548	0.665	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.20	278	4209	0.172	ng/ul#	96
26) Benzo(g,h,i)perylene	25.83	276	20151	0.721	ng/ul	97

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

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