

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012922.D
 Acq On : 06 Dec 2020 17:53
 Operator : CG/JU
 Sample : L4864-22
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B68

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:11:49 PM

Quant Time: Dec 07 03:31:40 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	2097	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	8673	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	4795	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	10183	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	7614	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	6274	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1605	0.12	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	3670	0.13	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.77	152	605	0.027	ng/ul#	79
9) Fluorene	15.12	166	1366	0.064	ng/ul	99
12) Phenanthrene	16.85	178	27045	0.792	ng/ul	97
13) Anthracene	16.95	178	19920	0.686	ng/ul	94
15) Fluoranthene	18.88	202	248972	6.137	ng/ul	99
17) Pyrene	19.26	202	241268	6.709	ng/ul	100
18) Benzo(a)anthracene	21.02	228	234588	7.447	ng/ul	98
19) Chrysene	21.07	228	259145	7.445	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	344330m	11.863	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	105640m	3.293	ng/ul	
23) Benzo(a)pyrene	23.05	252	194981	7.240	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.20	276	129961	3.674	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.20	278	33774	1.207	ng/ul	92
26) Benzo(g,h,i)perylene	25.83	276	102850	3.214	ng/ul	96

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

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