

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011755.D
 Acq On : 09 Sep 2020 02:45
 Operator : CG/JU
 Sample : L3843-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC1

Manual Integrations
 APPROVED

Sohil
 9/9/2020 3:46:36 PM

Quant Time: Sep 09 07:39:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 09 07:05:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2245	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8859	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	5213	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	10904m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	9181	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	9435	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4527	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	9104	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1952	0.074	ng/ul#	93
7) Acenaphthylene	14.09	152	1681	0.067	ng/ul#	89
12) Phenanthrene	17.16	178	1296	0.035	ng/ul#	89
13) Anthracene	17.25	178	1784	0.055	ng/ul#	90
15) Fluoranthene	19.18	202	3962	0.097	ng/ul	94
17) Pyrene	19.54	202	13514	0.331	ng/ul#	94
18) Benzo(a)anthracene	21.29	228	3397m	0.095	ng/ul	
19) Chrysene	21.34	228	6882	0.173	ng/ul	97
21) Benzo(b)fluoranthene	22.88	252	12290m	0.278	ng/ul	
22) Benzo(k)fluoranthene	22.92	252	4730m	0.104	ng/ul	
23) Benzo(a)pyrene	23.46	252	7744	0.203	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.81	276	4881	0.097	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.82	278	1439	0.035	ng/ul#	49
26) Benzo(g,h,i)perylene	26.50	276	4758	0.109	ng/ul#	88

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

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