

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022021\
 Data File : BN013689.D
 Acq On : 20 Feb 2021 13:33
 Operator : CG/JU
 Sample : M1412-24
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBGZ9

Manual Integrations
 APPROVED

mohammad
 2/22/2021 2:10:59 PM

Quant Time: Feb 22 06:08:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 16 16:31:21 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	4980	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	18736	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	12988	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	22312m	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	20823m	0.40	ng/ul	0.01
20) Perylene-d12	23.34	264	17353	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	7643	0.29	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	13200	0.24	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.23	166	1375	0.028	ng/ul#	91
11) Pentachlorophenol	16.59	266	112	0.041	ng/ul#	81
12) Phenanthrene	16.97	178	2493	0.038	ng/ul#	72
13) Anthracene	17.06	178	1194	0.022	ng/ul#	25
15) Fluoranthene	19.00	202	3385	0.047	ng/ul#	17
17) Pyrene	19.37	202	2182	0.029	ng/ul#	52
18) Benzo(a)anthracene	21.13	228	171788m	2.542	ng/ul	
19) Chrysene	21.17	228	277296m	3.771	ng/ul	
21) Benzo(b)fluoranthene	22.69	252	4766	0.067	ng/ul#	1
22) Benzo(k)fluoranthene	22.75	252	2651	0.036	ng/ul#	1
23) Benzo(a)pyrene	23.24	252	4485	0.079	ng/ul#	1
26) Benzo(g,h,i)perylene	26.16	276	1547	0.023	ng/ul#	1

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

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