

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112420\
 Data File : BN012727.D
 Acq On : 24 Nov 2020 21:39
 Operator : CG/JU
 Sample : L4751-23
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AY6

Manual Integrations
 APPROVED

mohammad
 11/25/2020 1:21:22 PM

Quant Time: Nov 25 03:21:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 25 01:45:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	887	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	3727	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2058	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	4094	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3203	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	3528	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	822	0.15	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	1798	0.17	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.89	178	1885	0.132	ng/ul	98
13) Anthracene	16.98	178	338	0.027	ng/ul#	66
15) Fluoranthene	18.92	202	5690	0.346	ng/ul	97
17) Pyrene	19.28	202	5576	0.389	ng/ul#	94
18) Benzo(a)anthracene	21.05	228	2355	0.205	ng/ul	95
19) Chrysene	21.10	228	3052	0.213	ng/ul	95
21) Benzo(b)fluoranthene	22.55	252	4269m	0.263	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	1563m	0.084	ng/ul	
23) Benzo(a)pyrene	23.08	252	2789	0.183	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.26	276	2113	0.106	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.26	278	458	0.029	ng/ul#	11
26) Benzo(g,h,i)perylene	25.89	276	2014	0.112	ng/ul#	87

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

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