

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027894.D
 Acq On : 21 Nov 2020 22:57
 Operator : JU/CG
 Sample : L4711-19
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B03

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:59:14 AM

Quant Time: Nov 23 05:00:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 04:04:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1155	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4689	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2949	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5360m	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2904	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2584	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2484	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4023	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	669	0.048	ng/ul#	79
5) 2-Methylnaphthalene	12.80	142	377	0.039	ng/ul	99
7) Acenaphthylene	14.68	152	688	0.048	ng/ul#	76
8) Acenaphthene	15.01	153	539	0.050	ng/ul	95
9) Fluorene	15.98	166	744	0.059	ng/ul#	97
11) Pentachlorophenol	17.31	266	39	0.021	ng/ul	88
12) Phenanthrene	17.70	178	21141	1.249	ng/ul	98
13) Anthracene	17.79	178	4112	0.252	ng/ul	97
15) Fluoranthene	19.68	202	72838	3.880	ng/ul	99
17) Pyrene	20.04	202	59179	4.013	ng/ul	98
18) Benzo(a)anthracene	21.75	228	28520	2.434	ng/ul	99
19) Chrysene	21.80	228	33553	2.940	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	40513m	3.731	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	14536m	1.367	ng/ul	
23) Benzo(a)pyrene	24.08	252	26339	2.774	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.68	276	18984	1.847	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.69	278	4757	0.556	ng/ul#	87
26) Benzo(g,h,i)perylene	27.46	276	18433	2.137	ng/ul	95

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

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