

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027923.D
 Acq On : 23 Nov 2020 21:02
 Operator : JU/CG
 Sample : L4711-17DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW9DL

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:45:06 PM

Quant Time: Nov 24 01:10:03 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 : Tue Nov 24 00:42:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1237	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4913	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2810	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5085	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2941	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2500	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	425	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	700	0.05	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	15.00	153	250	0.024	ng/ul#	93
9) Fluorene	15.98	166	292	0.024	ng/ul#	88
12) Phenanthrene	17.70	178	8496	0.529	ng/ul	98
13) Anthracene	17.79	178	2217	0.143	ng/ul	94
15) Fluoranthene	19.68	202	17389	0.976	ng/ul	98
17) Pyrene	20.03	202	14469	0.969	ng/ul	97
18) Benzo(a)anthracene	21.75	228	6303	0.531	ng/ul	98
19) Chrysene	21.80	228	6141	0.531	ng/ul	97
21) Benzo(b)fluoranthene	23.45	252	6534m	0.622	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	2107m	0.205	ng/ul	
23) Benzo(a)pyrene	24.08	252	4482	0.488	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.67	276	2703	0.272	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.68	278	722	0.087	ng/ul#	40
26) Benzo(g,h,i)perylene	27.45	276	2528	0.303	ng/ul	95

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

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