

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027967.D
 Acq On : 25 Nov 2020 17:11
 Operator : JU/CG
 Sample : L4751-13DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B85DL

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:43:44 AM

Quant Time: Nov 25 17:43:45 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 : Wed Nov 25 10:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	1021	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4118	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2315	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4184	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2413	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2286	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	153	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	227	0.02	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	15.00	153	381	0.045	ng/ul	95
9) Fluorene	15.97	166	398	0.041	ng/ul#	93
12) Phenanthrene	17.69	178	13192	0.998	ng/ul	98
13) Anthracene	17.79	178	2085	0.164	ng/ul	94
15) Fluoranthene	19.67	202	45750	3.122	ng/ul	98
17) Pyrene	20.03	202	36919	3.013	ng/ul	97
18) Benzo(a)anthracene	21.74	228	18392	1.889	ng/ul	100
19) Chrysene	21.80	228	24134	2.545	ng/ul	98
21) Benzo(b)fluoranthene	23.44	252	31482m	3.277	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	11991m	1.275	ng/ul	
23) Benzo(a)pyrene	24.07	252	17766	2.115	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.67	276	13754	1.512	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.67	278	3540	0.467	ng/ul	97
26) Benzo(g,h,i)perylene	27.44	276	12139	1.591	ng/ul	94

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

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