

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027865.D
 Acq On : 20 Nov 2020 22:50
 Operator : JU/CG
 Sample : L4710-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BF0

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:57:48 AM

Quant Time: Nov 21 03:10:20 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 : Sat Nov 21 02:40:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	950	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3900	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2449	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	4800m	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2911	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2261	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1277	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	2739	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	328	0.028	ng/ul#	61
5) 2-Methylnaphthalene	12.80	142	190	0.024	ng/ul	96
7) Acenaphthylene	14.68	152	277	0.023	ng/ul#	56
12) Phenanthrene	17.70	178	4796	0.316	ng/ul	99
13) Anthracene	17.79	178	883	0.060	ng/ul#	82
15) Fluoranthene	19.68	202	15778	0.938	ng/ul	98
17) Pyrene	20.04	202	14217	0.962	ng/ul	98
18) Benzo(a)anthracene	21.75	228	6229	0.530	ng/ul	97
19) Chrysene	21.80	228	8038	0.703	ng/ul	98
21) Benzo(b)fluoranthene	23.45	252	9512m	1.001	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	3269m	0.351	ng/ul	
23) Benzo(a)pyrene	24.08	252	5654	0.681	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.68	276	4018	0.447	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.68	278	1048	0.140	ng/ul#	42
26) Benzo(g,h,i)perylene	27.45	276	3833	0.508	ng/ul	97

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

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