

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027899.D
 Acq On : 22 Nov 2020 03:12
 Operator : JU/CG
 Sample : L4711-24
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BF6

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:59:27 AM

Quant Time: Nov 23 05:52:31 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	1086	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4238	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2547	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4266m	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2278	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2072	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1743	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	2601	0.23	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.22	128	857	0.067	ng/ul#	85
5) 2-Methylnaphthalene	12.80	142	904	0.105	ng/ul	100
7) Acenaphthylene	14.68	152	954	0.077	ng/ul#	83
8) Acenaphthene	15.01	153	6415	0.689	ng/ul	99
9) Fluorene	15.98	166	6793	0.629	ng/ul	100
12) Phenanthrene	17.70	178	120154	8.916	ng/ul	97
13) Anthracene	17.79	178	26223	2.017	ng/ul	98
15) Fluoranthene	19.68	202	168278	11.262	ng/ul	100
17) Pyrene	20.04	202	129901	11.228	ng/ul	99
18) Benzo(a)anthracene	21.74	228	55631	6.053	ng/ul	100
19) Chrysene	21.80	228	57326	6.403	ng/ul	98
21) Benzo(b)fluoranthene	23.45	252	67128m	7.709	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	21484m	2.519	ng/ul	
23) Benzo(a)pyrene	24.08	252	43841	5.758	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.68	276	31010	3.762	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.68	278	7762	1.131	ng/ul#	92
26) Benzo(a,h,i)perylene	27.45	276	30241	4.373	ng/ul	93

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

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