

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027881.D
 Acq On : 21 Nov 2020 14:58
 Operator : JU/CG
 Sample : L4854-06
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD4

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:58:42 AM

Quant Time: Nov 23 03:52:37 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 02:58:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1175	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	4800	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2926	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5586	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	2969	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2597	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1107	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	1956	0.13	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	16406	1.140	ng/ul	97
5) 2-Methylnaphthalene	12.80	142	30952	3.162	ng/ul	99
7) Acenaphthylene	14.67	152	1227	0.086	ng/ul#	41
8) Acenaphthene	15.02	153	1272	0.119	ng/ul#	59
9) Fluorene	15.98	166	1796	0.145	ng/ul#	82
12) Phenanthrene	17.70	178	34737	1.969	ng/ul	99
13) Anthracene	17.79	178	2920m	0.172	ng/ul	
15) Fluoranthene	19.68	202	14194	0.725	ng/ul	98
17) Pyrene	20.04	202	17017	1.129	ng/ul	97
18) Benzo(a)anthracene	21.75	228	5552	0.463	ng/ul#	60
19) Chrysene	21.80	228	8938	0.766	ng/ul#	70
21) Benzo(b)fluoranthene	23.45	252	7004m	0.642	ng/ul	
22) Benzo(k)fluoranthene	23.50	252	2455m	0.230	ng/ul	
23) Benzo(a)pyrene	24.08	252	4450	0.466	ng/ul#	43
24) Indeno(1,2,3-cd)pyrene	26.68	276	3441	0.333	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.69	278	1002	0.116	ng/ul#	1
26) Benzo(a,h,i)perylene	27.46	276	5233	0.604	ng/ul#	79

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

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