

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112520\
 Data File : BM027969.D
 Acq On : 25 Nov 2020 18:24
 Operator : JU/CG
 Sample : L4711-01DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B36DL

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:43:49 AM

Quant Time: Nov 26 00:55:24 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	1022	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4039	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2284	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4040	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	2277	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	327	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	457	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	443	0.037	ng/ul#	70
5) 2-Methylnaphthalene	12.79	142	311	0.038	ng/ul	100
7) Acenaphthylene	14.66	152	523	0.047	ng/ul#	76
8) Acenaphthene	15.00	153	2337	0.280	ng/ul	99
9) Fluorene	15.97	166	2871	0.296	ng/ul	99
12) Phenanthrene	17.70	178	57654	4.518	ng/ul	96
13) Anthracene	17.78	178	6898	0.560	ng/ul	99
15) Fluoranthene	19.67	202	89677	6.337	ng/ul	100
17) Pyrene	20.03	202	67740	5.528	ng/ul	97
18) Benzo(a)anthracene	21.74	228	28681	2.946	ng/ul	100
19) Chrysene	21.80	228	33413	3.523	ng/ul	98
21) Benzo(b)fluoranthene	23.44	252	39315m	4.108	ng/ul	
22) Benzo(k)fluoranthene	23.48	252	14652m	1.564	ng/ul	
23) Benzo(a)pyrene	24.07	252	25744	3.077	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.67	276	17890	1.975	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.67	278	4398	0.583	ng/ul	97
26) Benzo(a,h,i)perylene	27.44	276	16985	2.235	ng/ul	94

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

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