

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025304.D
 Acq On : 06 Mar 2020 02:10
 Operator : CG/JU
 Sample : L1568-04
 Misc :
 ALS Vial : 101 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDA7

Manual Integrations
 APPROVED

mohammad
 3/6/2020 10:15:29 PM

Quant Time: Mar 06 03:40:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3576	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	15743	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11780	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	28166m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	27514	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	25861	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	7030	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	22560	0.25	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	1060	0.022	ng/ul	98
8) Acenaphthene	14.36	153	1241	0.029	ng/ul	97
9) Fluorene	15.35	166	1229	0.023	ng/ul	95
12) Phenanthrene	17.08	178	7269	0.080	ng/ul	99
13) Anthracene	17.17	178	2585	0.032	ng/ul	97
15) Fluoranthene	19.10	202	12675	0.112	ng/ul	98
17) Pyrene	19.46	202	14104m	0.135	ng/ul	
18) Benzo(a)anthracene	21.21	228	7793	0.077	ng/ul	97
19) Chrysene	21.26	228	9584	0.086	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	16740m	0.162	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	5766m	0.053	ng/ul	
23) Benzo(a)pyrene	23.32	252	6460	0.074	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.59	276	3872	0.036	ng/ul#	100
26) Benzo(g,h,i)perylene	26.26	276	3515	0.038	ng/ul#	93

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

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