

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025302.D
 Acq On : 06 Mar 2020 00:58
 Operator : CG/JU
 Sample : L1568-06
 Misc :
 ALS Vial : 99 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDA9

Manual Integrations
 APPROVED

mohammad
 3/6/2020 10:15:23 PM

Quant Time: Mar 06 03:34:14 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	4052	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	17884	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	13461	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	32086m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	28692	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	24961	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	7136	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	23863	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	1762	0.033	ng/ul	94
8) Acenaphthene	14.36	153	6330	0.129	ng/ul	98
9) Fluorene	15.35	166	1815	0.030	ng/ul	95
12) Phenanthrene	17.08	178	3057	0.030	ng/ul	98
13) Anthracene	17.17	178	2653	0.029	ng/ul#	96
15) Fluoranthene	19.10	202	23013	0.178	ng/ul	97
17) Pyrene	19.46	202	40599m	0.373	ng/ul	
18) Benzo(a)anthracene	21.21	228	18039	0.171	ng/ul	98
19) Chrysene	21.26	228	34722	0.300	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	40137m	0.402	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	14325m	0.137	ng/ul	
23) Benzo(a)pyrene	23.32	252	16624	0.198	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.59	276	8636	0.083	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	2375	0.027	ng/ul#	66
26) Benzo(g,h,i)perylene	26.25	276	4801	0.054	ng/ul	95

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

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