

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025303.D
 Acq On : 06 Mar 2020 01:34
 Operator : CG/JU
 Sample : L1568-12
 Misc :
 ALS Vial : 100 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDB3

Manual Integrations
 APPROVED

mohammad
 3/6/2020 10:15:26 PM

Quant Time: Mar 06 03:37:40 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	4290	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	18919	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	14110	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	33384m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	32274	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	28607	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	7282	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	25252	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	1380	0.024	ng/ul	95
11) Pentachlorophenol	16.70	266	198	0.023	ng/ul	90
13) Anthracene	17.17	178	2399	0.025	ng/ul#	92
15) Fluoranthene	19.10	202	13257	0.098	ng/ul	99
17) Pyrene	19.46	202	12649	0.103	ng/ul	99
18) Benzo(a)anthracene	21.21	228	8819	0.074	ng/ul	98
19) Chrysene	21.26	228	13648m	0.105	ng/ul	
21) Benzo(b)fluoranthene	22.76	252	21092m	0.184	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	5116m	0.043	ng/ul	
23) Benzo(a)pyrene	23.32	252	7969	0.083	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.59	276	3925	0.033	ng/ul#	100
26) Benzo(g,h,i)perylene	26.25	276	2526	0.025	ng/ul	92

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

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