

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025382.D
 Acq On : 08 Mar 2020 12:36
 Operator : CG/JU
 Sample : L1616-18
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDR2

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:03:20 PM

Quant Time: Mar 09 02:00:10 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	7.66	152	4318	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.43	136	18152	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.29	164	13084	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.04	188	30331m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	29581	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	29355	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	9689	0.29	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	28410	0.29	ng/ul	-0.01
Target Compounds				Ovalue		
12) Phenanthrene	17.08	178	2124	0.022	ng/ul	95
13) Anthracene	17.17	178	2186	0.025	ng/ul#	96
15) Fluoranthene	19.09	202	4705	0.038	ng/ul	98
17) Pyrene	19.46	202	5787	0.052	ng/ul#	96
19) Chrysene	21.26	228	3835	0.032	ng/ul	97
21) Benzo(b)fluoranthene	22.76	252	7723m	0.066	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.58	276	3063	0.025	ng/ul#	100
26) Benzo(g,h,i)perylene	26.24	276	2444	0.023	ng/ul#	91

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

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