

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025331.D
 Acq On : 07 Mar 2020 03:18
 Operator : CG/JU
 Sample : L1612-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDL1

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:00:39 PM

Quant Time: Mar 07 05:20:05 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 : Sat Mar 07 00:01:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3265	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14722	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10985	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	25550m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	24546	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	22493	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	9250	0.34	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	25627	0.31	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.36	153	3036	0.076	ng/ul	100
9) Fluorene	15.35	166	6356	0.127	ng/ul	98
12) Phenanthrene	17.08	178	40705	0.494	ng/ul	95
15) Fluoranthene	19.10	202	23691	0.230	ng/ul	96
17) Pyrene	19.46	202	14490m	0.156	ng/ul	
18) Benzo(a)anthracene	21.21	228	2175	0.024	ng/ul	97
19) Chrysene	21.26	228	2433	0.025	ng/ul	98

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

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