

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025351.D
 Acq On : 07 Mar 2020 15:52
 Operator : CG/JU
 Sample : L1616-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDP7

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:01:42 PM

Quant Time: Mar 07 21:43:25 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 20:27:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4651	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	21074	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.30	164	15526	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	35688m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	33095	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	29214	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	13549	0.34	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	38682	0.33	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.02	152	1680	0.023	ng/ul#	92
17) Pyrene	19.46	202	11809	0.094	ng/ul	98
18) Benzo(a)anthracene	21.21	228	16173	0.133	ng/ul	100
19) Chrysene	21.26	228	14270m	0.107	ng/ul	
21) Benzo(b)fluoranthene	22.76	252	28234m	0.242	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	10764m	0.088	ng/ul	
23) Benzo(a)pyrene	23.32	252	13564	0.138	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.58	276	5486	0.045	ng/ul#	100
26) Benzo(g,h,i)perylene	26.25	276	4125	0.040	ng/ul	94

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

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