

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025519.D
 Acq On : 12 Mar 2020 12:53
 Operator : CG/JU
 Sample : L1677-20
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CB7

Manual Integrations
 APPROVED

Quant Time: Mar 13 04:48:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2549	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	10893	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	7664	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	16265	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	10874	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	11140	0.40	ng/ul	0.00

mohammad

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5829	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	12947	0.29	ng/ul	0.00

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Target Compounds						Ovalue
12) Phenanthrene	17.07	178	4732	0.092	ng/ul	95
15) Fluoranthene	19.09	202	10881	0.195	ng/ul	97
17) Pyrene	19.45	202	8983	0.184	ng/ul	98
18) Benzo(a)anthracene	21.20	228	3020m	0.078	ng/ul	
19) Chrysene	21.25	228	4365	0.100	ng/ul	97
21) Benzo(b)fluoranthene	22.75	252	4737m	0.107	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	2355m	0.051	ng/ul	
23) Benzo(a)pyrene	23.30	252	3092	0.084	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.57	276	2580	0.052	ng/ul#	63
26) Benzo(g,h,i)perylene	26.23	276	2488	0.058	ng/ul#	83

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

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