

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100520\
 Data File : BM027613.D
 Acq On : 05 Oct 2020 19:49
 Operator : JU/CG
 Sample : L4184-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ6

Manual Integrations
 APPROVED

mohammad
 10/6/2020 3:23:54 PM

Quant Time: Oct 06 01:23:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 06 01:23:11 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	1234	0.40	ng/ul	0.00
2) Naphthalene-d8	10.579	136	3575	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.423	164	2557	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.168	188	6191m	0.40	ng/ul	0.00
16) Chrysene-d12	21.345	240	4652	0.40	ng/ul	0.00
20) Perylene-d12	23.605	264	7279	0.40	ng/ul	# 0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.167	152	1299	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.190	212	3976	0.23	ng/ul	0.00
Target Compounds						
						Qvalue
12) Phenanthrene	17.209	178	879	0.05	ng/ul#	85
15) Fluoranthene	19.220	202	2759	0.12	ng/ul	92
17) Pyrene	19.583	202	2715m	0.13	ng/ul	
18) Benzo(a)anthracene	21.328	228	931	0.05	ng/ul#	88
19) Chrysene	21.381	228	1154	0.06	ng/ul#	94
21) Benzo(b)fluoranthene	22.926	252	1719m	0.06	ng/ul	
23) Benzo(a)pyrene	23.505	252	1515	0.06	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	25.880	276	1200	0.04	ng/ul#	94
26) Benzo(g,h,i)perylene	26.578	276	1166	0.04	ng/ul#	73

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

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