

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100220\
 Data File : BM027595.D
 Acq On : 02 Oct 2020 23:36
 Operator : JU/CG
 Sample : L4166-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

mohammad
 10/5/2020 10:29:23 AM

Quant Time: Oct 03 01:38:41 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 : Fri Oct 02 13:04:03 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	1324	0.40	ng/ul	0.00
2) Naphthalene-d8	10.581	136	3831	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.429	164	2823	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.169	188	6934m	0.40	ng/ul	0.00
16) Chrysene-d12	21.343	240	7463	0.40	ng/ul	0.00
20) Perylene-d12	23.605	264	8198	0.40	ng/ul	# 0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.169	152	1368	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.192	212	4218	0.22	ng/ul	0.00
Target Compounds						
						Qvalue
15) Fluoranthene	19.226	202	690	0.03	ng/ul	77
17) Pyrene	19.584	202	833	0.02	ng/ul#	74
18) Benzo(a)anthracene	21.326	228	883	0.03	ng/ul#	83
19) Chrysene	21.379	228	1635	0.05	ng/ul#	90
21) Benzo(b)fluoranthene	22.927	252	2452m	0.07	ng/ul	
22) Benzo(k)fluoranthene	22.968	252	912m	0.03	ng/ul	
23) Benzo(a)pyrene	23.506	252	1592	0.05	ng/ul#	36
24) Indeno(1,2,3-cd)pyrene	25.882	276	1056	0.03	ng/ul#	93
26) Benzo(g,h,i)perylene	26.573	276	959	0.03	ng/ul#	68

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

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