

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092721\
 Data File : BM032238.D
 Acq On : 28 Sep 2021 12:44
 Operator : CG/JU
 Sample : M3739-07
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AE2

Manual Integrations
 APPROVED

mohammad
 9/29/2021 12:52:12 PM

Quant Time: Sep 28 13:26:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 27 13:37:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	7025	0.400	ng/ul	0.00
4) Naphthalene-d8	10.451	136	21800	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.303	164	10839	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.028	188	19421	0.400	ng/ul	0.00
17) Chrysene-d12	21.192	240	14846	0.400	ng/ul	# 0.00
23) Perylene-d12	23.357	264	15312	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.982	96	39287	5.202	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.038	152	9807	0.338	ng/ul	0.00
18) Fluoranthene-d10	19.049	212	19463	0.483	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.070	178	1748	0.028	ng/ul#	80
19) Fluoranthene	19.079	202	5404	0.087	ng/ul#	94
20) Pyrene	19.441	202	7023	0.113	ng/ul#	96
21) Benzo(a)anthracene	21.175	228	2130	0.041	ng/ul#	83
22) Chrysene	21.229	228	2301	0.037	ng/ul#	83
24) Benzo(b)fluoranthene	22.713	252	2315m	0.029	ng/ul	
26) Benzo(a)pyrene	23.268	252	3839	0.064	ng/ul#	1

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

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