

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015504.D
 Acq On : 12 Jul 2021 23:49
 Operator : CG/JU
 Sample : M2854-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD4

Manual Integrations
 APPROVED

mohammad
 7/13/2021 1:44:53 PM

Quant Time: Jul 13 04:58:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 12 11:02:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	3802	0.400	ng/ul	0.00
4) Naphthalene-d8	10.595	136	12460	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.442	164	6144	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.192	188	11436	0.400	ng/ul	0.00
17) Chrysene-d12	21.383	240	9192	0.400	ng/ul	# 0.00
23) Perylene-d12	23.739	264	9061	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	7895	1.842	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	3131	0.160	ng/ul	0.00
18) Fluoranthene-d10	19.226	212	5864	0.194	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.234	178	1155	0.028	ng/ul#	87
19) Fluoranthene	19.253	202	1998	0.047	ng/ul#	88
20) Pyrene	19.616	202	1533m	0.036	ng/ul	
21) Benzo(a)anthracene	21.368	228	1068	0.031	ng/ul#	78
22) Chrysene	21.421	228	897	0.022	ng/ul#	86
24) Benzo(b)fluoranthene	23.031	252	879m	0.021	ng/ul	

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

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