

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071521\
 Data File : BM030981.D
 Acq On : 15 Jul 2021 15:04
 Operator : CG/JU
 Sample : M2910-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW2

Manual Integrations
 APPROVED

mohammad
 7/16/2021 10:56:31 AM

Quant Time: Jul 15 15:36:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM071321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 15 13:02:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.656	152	988	0.400	ng/ul	0.00
4) Naphthalene-d8	10.423	136	3287	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.284	164	2330	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.030	188	5052m	0.400	ng/ul	0.00
17) Chrysene-d12	21.219	240	4880	0.400	ng/ul	0.00
23) Perylene-d12	23.408	264	5531	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.227	96	2251	2.216	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.015	152	1634	0.286	ng/ul	0.00
18) Fluoranthene-d10	19.061	212	5060	0.308	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.262	88	167	0.147	ng/ul	89
5) Naphthalene	10.472	128	348	0.035	ng/ul#	52
11) Acenaphthene	14.348	153	300	0.036	ng/ul#	87
12) Fluorene	15.338	166	636	0.065	ng/ul#	94
14) Pentachlorophenol	16.683	266	139887	78.385	ng/ul	99
15) Phenanthrene	17.071	178	380	0.024	ng/ul#	40

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

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