

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030969.D
 Acq On : 14 Jul 2021 20:37
 Operator : CG/JU
 Sample : M2910-03
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW3

Manual Integrations
 APPROVED

mohammad
 7/15/2021 11:31:10 AM

Quant Time: Jul 15 04:07:51 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 : Tue Jul 13 14:01:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	1128	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.428	136	3933	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.292	164	2945	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.037	188	6202m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.224	240	5938	0.400	ng/ul	-0.01
23) Perylene-d12	23.413	264	6053	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	5436	4.686	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.020	152	2511	0.367	ng/ul	-0.01
18) Fluoranthene-d10	19.066	212	8895	0.445	ng/ul	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.269	88	305	0.236	ng/ul#	41
11) Acenaphthene	14.353	153	258	0.025	ng/ul#	91
14) Pentachlorophenol	16.690	266	215427	98.330	ng/ul	99
15) Phenanthrene	17.079	178	521	0.027	ng/ul#	36
16) Anthracene	17.172	178	553	0.030	ng/ul#	50

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

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