

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030968.D
 Acq On : 14 Jul 2021 20:00
 Operator : CG/JU
 Sample : M2910-01
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW2

Manual Integrations
 APPROVED

mohammad
 7/15/2021 11:31:07 AM

Quant Time: Jul 15 04:07:42 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	1332	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.427	136	4564	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.291	164	3221	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.037	188	6541m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.225	240	5983	0.400	ng/ul	-0.01
23) Perylene-d12	23.417	264	6045	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	6440	4.702	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.019	152	2486	0.313	ng/ul	-0.01
18) Fluoranthene-d10	19.069	212	7222	0.359	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.268	88	262	0.171	ng/ul#	71
5) Naphthalene	10.476	128	483	0.035	ng/ul#	56
11) Acenaphthene	14.356	153	437	0.038	ng/ul#	91
12) Fluorene	15.342	166	875	0.065	ng/ul#	95
14) Pentachlorophenol	16.686	266	204700	88.591	ng/ul	99
15) Phenanthrene	17.078	178	530	0.026	ng/ul#	50

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

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