

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030959.D
 Acq On : 14 Jul 2021 12:57
 Operator : CG/JU
 Sample : M2905-23
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBK46

Manual Integrations
 APPROVED

mohammad
 7/14/2021 4:35:49 PM

Quant Time: Jul 14 13:27:16 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	1078	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.433	136	3631	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.296	164	2582	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.041	188	5417m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.227	240	4857	0.400	ng/ul	0.00
23) Perylene-d12	23.421	264	5416	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.237	96	4987	4.499	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.025	152	2303	0.364	ng/ul	0.00
18) Fluoranthene-d10	19.070	212	6786	0.415	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.690	266	93	0.049	ng/ul#	90
15) Phenanthrene	17.082	178	676	0.040	ng/ul#	77
16) Anthracene	17.173	178	336	0.021	ng/ul#	53
19) Fluoranthene	19.100	202	626	0.029	ng/ul#	61
20) Pyrene	19.462	202	592	0.027	ng/ul#	53
21) Benzo(a)anthracene	21.212	228	591	0.028	ng/ul#	85
22) Chrysene	21.263	228	540	0.025	ng/ul#	83
24) Benzo(b)fluoranthene	22.767	252	773	0.030	ng/ul#	31
25) Benzo(k)fluoranthene	22.805	252	748	0.028	ng/ul#	5
26) Benzo(a)pyrene	23.324	252	586m	0.026	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.599	276	979	0.032	ng/ul#	69
28) Dibenzo(a,h)anthracene	25.613	278	655	0.026	ng/ul#	30
29) Benzo(g,h,i)perylene	26.263	276	790	0.030	ng/ul#	60

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

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