

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092721\
 Data File : BM032221.D
 Acq On : 28 Sep 2021 00:57
 Operator : CG/JU
 Sample : M3919-07
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GB6Y6

Quant Time: Sep 28 02:31:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 27 13:37:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.660	152	7456	0.400	ng/ul	0.00
4) Naphthalene-d8	10.451	136	23198	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.304	164	11511	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.033	188	20733	0.400	ng/ul	0.00
17) Chrysene-d12	21.191	240	15569	0.400	ng/ul	0.00
23) Perylene-d12	23.356	264	15255	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.989	96	42013	5.241	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.044	152	10570	0.342	ng/ul	0.00
18) Fluoranthene-d10	19.049	212	19530	0.462	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.027	88	645	0.078	ng/ul#	74
12) Fluorene	15.347	166	1354	0.028	ng/ul#	95
14) Pentachlorophenol	16.700	266	271	0.052	ng/ul	98
15) Phenanthrene	17.071	178	3653	0.054	ng/ul#	93
16) Anthracene	17.161	178	3733	0.067	ng/ul#	93
19) Fluoranthene	19.080	202	4965	0.076	ng/ul	95
20) Pyrene	19.442	202	5259	0.081	ng/ul#	92
21) Benzo(a)anthracene	21.174	228	3786	0.069	ng/ul	95
22) Chrysene	21.227	228	4270	0.066	ng/ul	98
24) Benzo(b)fluoranthene	22.709	252	4272	0.054	ng/ul#	41
25) Benzo(k)fluoranthene	22.753	252	3726	0.049	ng/ul#	37
26) Benzo(a)pyrene	23.259	252	3408	0.057	ng/ul#	36
27) Indeno(1,2,3-cd)pyrene	25.498	276	3767	0.050	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.498	278	3033	0.049	ng/ul#	61
29) Benzo(g,h,i)perylene	26.157	276	3570	0.050	ng/ul#	81

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

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