

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031380.D
 Acq On : 02 Aug 2021 17:56
 Operator : CG/JU
 Sample : M3123-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0074

Quant Time: Aug 02 18:27:03 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 02 11:29:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.601	152	1307	0.40	ng/ul	0.00
4) Naphthalene-d8	10.365	136	4990	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.231	164	2885	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.978	188	5261	0.40	ng/ul	0.00
17) Chrysene-d12	21.171	240	3939	0.40	ng/ul	0.00
23) Perylene-d12	23.334	264	4005	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	2088	1.44	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.957	152	1063	0.13	ng/ul	0.00
18) Fluoranthene-d10	19.012	212	1942	0.15	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.410	128	1580	0.11	ng/ul#	93
7) 2-Methylnaphthalene	12.034	142	1034	0.10	ng/ul	100
8) 1-Methylnaphthalene	12.254	142	741	0.07	ng/ul	97
14) Pentachlorophenol	16.638	266	77	0.05	ng/ul	93
15) Phenanthrene	17.020	178	4693	0.28	ng/ul	98
16) Anthracene	17.110	178	1004	0.06	ng/ul#	86
19) Fluoranthene	19.043	202	2339	0.14	ng/ul#	92
20) Pyrene	19.405	202	3278	0.19	ng/ul#	94
21) Benzo(a)anthracene	21.156	228	2693	0.17	ng/ul#	86
22) Chrysene	21.207	228	5066	0.31	ng/ul#	93
24) Benzo(b)fluoranthene	22.689	252	2919	0.16	ng/ul#	75
25) Benzo(k)fluoranthene	22.733	252	1245	0.07	ng/ul#	57
26) Benzo(a)pyrene	23.239	252	2019	0.13	ng/ul#	34
27) Indeno(1,2,3-cd)pyrene	25.473	276	1674	0.08	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.490	278	1558	0.09	ng/ul#	61
29) Benzo(g,h,i)perylene	26.127	276	1267	0.07	ng/ul#	74

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

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