

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031266.D
 Acq On : 30 Jul 2021 01:57
 Operator : CG/JU
 Sample : M3084-14
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0032

Quant Time: Jul 30 03:43:01 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 : Wed Jul 28 15:20:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.614	152	1548	0.400	ng/ul	0.00
4) Naphthalene-d8	10.374	136	5752	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.242	164	3463	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.988	188	6620	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.178	240	4994	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	5187	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	4561	2.653	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.971	152	1614	0.167	ng/ul	0.00
18) Fluoranthene-d10	19.020	212	3313	0.199	ng/ul	0.00
Target Compounds						
					Qvalue	
7) 2-Methylnaphthalene	12.043	142	318	0.027	ng/ul	98
15) Phenanthrene	17.030	178	1233	0.059	ng/ul#	90
19) Fluoranthene	19.050	202	2080	0.098	ng/ul#	87
20) Pyrene	19.412	202	2025	0.094	ng/ul#	84
21) Benzo(a)anthracene	21.161	228	1334	0.066	ng/ul#	89
22) Chrysene	21.214	228	1551	0.074	ng/ul#	90
24) Benzo(b)fluoranthene	22.701	252	1839	0.079	ng/ul#	46
25) Benzo(k)fluoranthene	22.740	252	815	0.034	ng/ul#	29
26) Benzo(a)pyrene	23.246	252	1185	0.058	ng/ul#	25
27) Indeno(1,2,3-cd)pyrene	25.487	276	1073	0.041	ng/ul#	73
28) Dibenzo(a,h)anthracene	25.502	278	581	0.027	ng/ul#	12
29) Benzo(g,h,i)perylene	26.141	276	1006	0.045	ng/ul#	67

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

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