

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031309.D
 Acq On : 31 Jul 2021 13:16
 Operator : CG/JU
 Sample : M3091-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0036

Quant Time: Aug 02 01:08:06 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 : Fri Jul 30 14:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	1792	0.400	ng/ul	0.00
4) Naphthalene-d8	10.370	136	6597	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.239	164	4005	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.985	188	7799	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.176	240	6166	0.400	ng/ul	# 0.00
23) Perylene-d12	23.341	264	6270	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	7101	3.568	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.966	152	2491	0.224	ng/ul	0.00
18) Fluoranthene-d10	19.016	212	5290	0.257	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.027	178	682	0.028	ng/ul#	81
16) Anthracene	17.120	178	649	0.028	ng/ul#	74
19) Fluoranthene	19.046	202	1048	0.040	ng/ul#	72
20) Pyrene	19.412	202	1101	0.041	ng/ul#	75
21) Benzo(a)anthracene	21.159	228	1510	0.061	ng/ul#	90
22) Chrysene	21.212	228	1806	0.070	ng/ul#	92
24) Benzo(b)fluoranthene	22.696	252	1389	0.049	ng/ul#	33
25) Benzo(k)fluoranthene	22.737	252	1606	0.055	ng/ul#	45
26) Benzo(a)pyrene	23.246	252	1363	0.055	ng/ul#	28
27) Indeno(1,2,3-cd)pyrene	25.492	276	1421	0.044	ng/ul#	64
28) Dibenzo(a,h)anthracene	25.499	278	1529	0.059	ng/ul#	63
29) Benzo(g,h,i)perylene	26.137	276	1484	0.054	ng/ul#	74

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

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