

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090923\
 Data File : BM041757.D
 Acq On : 09 Sep 2023 19:56
 Operator : MA/JU
 Sample : 04247-08
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBJ81

Quant Time: Sep 09 23:34:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM090523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 09 23:26:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.000	152	58	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.812	136	89	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.634	164	29	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.383	188	59	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.556	240	22	0.400	ng/ul	# 0.00
23) Perylene-d12	23.974	264	9	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.427	96	407	4.592	ng/ul	0.03
6) 2-Methylnaphthalene-d10	12.390	152	38	0.332	ng/ul	0.00
18) Fluoranthene-d10	19.399	212	32	0.300	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.465	88	3011	32.699	ng/ul#	86
8) 1-Methylnaphthalene	12.682	142	7	0.042	ng/ul#	48
10) Acenaphthylene	14.356	152	4	0.022	ng/ul#	1
11) Acenaphthene	14.698	153	4	0.029	ng/ul#	76
12) Fluorene	15.670	166	18	0.115	ng/ul#	53
14) Pentachlorophenol	17.008	266	2	0.085	ng/ul#	51
15) Phenanthrene	17.413	178	18	0.072	ng/ul#	1
16) Anthracene	17.510	178	41	0.192	ng/ul#	1
19) Fluoranthene	19.431	202	39	0.150	ng/ul#	1
20) Pyrene	19.798	202	58	0.225	ng/ul#	1
21) Benzo(a)anthracene	21.536	228	13	0.072	ng/ul#	1
22) Chrysene	21.585	228	11	0.057	ng/ul#	1
24) Benzo(b)fluoranthene	23.216	252	15	0.186	ng/ul#	1
25) Benzo(k)fluoranthene	23.278	252	30	0.381	ng/ul#	1
26) Benzo(a)pyrene	23.863	252	21	0.321	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.485	276	4	0.045	ng/ul#	70
28) Dibenzo(a,h)anthracene	26.508	278	8	0.127	ng/ul#	1
29) Benzo(g,h,i)perylene	27.269	276	13	0.163	ng/ul#	1

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

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