

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100423\
 Data File : BM042203.D
 Acq On : 05 Oct 2023 01:32
 Operator : MA/JU
 Sample : 04588-22
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBJZ9

Quant Time: Oct 05 03:00:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	118	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.741	136	342	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.574	164	366	0.400	ng/ul	#-0.01
13) Phenanthrene-d10	17.329	188	677	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.507	240	506	0.400	ng/ul	# 0.00
23) Perylene-d12	0.000	264	0	0.000	ng/ul	-23.92
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.355	96	37460	238.387	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.324	152	4585	9.775	ng/ul	-0.01
18) Fluoranthene-d10	19.352	212	7886	4.151	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.393	88	2565	15.631	ng/ul#	89
5) Naphthalene	10.796	128	953	0.835	ng/ul#	45
7) 2-Methylnaphthalene	12.401	142	181	0.265	ng/ul#	78
8) 1-Methylnaphthalene	12.621	142	185	0.269	ng/ul	94
10) Acenaphthylene	14.277	152	130	0.056	ng/ul#	1
11) Acenaphthene	14.638	153	37	0.021	ng/ul#	65
12) Fluorene	15.568	166	2049	1.039	ng/ul#	16
14) Pentachlorophenol	16.966	266	33	0.108	ng/ul#	82
15) Phenanthrene	17.371	178	157	0.056	ng/ul#	10
16) Anthracene	17.464	178	189	0.074	ng/ul#	15
19) Fluoranthene	19.385	202	235	0.061	ng/ul#	22
20) Pyrene	19.719	202	2073	0.507	ng/ul#	1
21) Benzo(a)anthracene	21.492	228	322	0.099	ng/ul#	58
22) Chrysene	21.545	228	342	0.108	ng/ul#	61

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

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