

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037677.D
 Acq On : 23 Nov 2022 13:10
 Operator : CG/JU
 Sample : N5568-09
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3D68

Quant Time: Nov 24 01:49:22 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 22:36:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.125	152	6314	0.400	ng/ul	0.00
4) Naphthalene-d8	10.942	136	18465	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.743	164	10614	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.475	188	24098	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.636	240	17554	0.400	ng/ul	# 0.00
23) Perylene-d12	24.127	264	16393	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	3008	0.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.520	152	8013	0.262	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	19749	0.340	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.991	128	31288	0.558	ng/ul	97
7) 2-Methylnaphthalene	12.592	142	20661	0.542	ng/ul	100
12) Fluorene	15.784	166	31512	0.636	ng/ul	99
14) Pentachlorophenol	17.120	266	7897	1.460	ng/ul	99
15) Phenanthrene	17.517	178	56193	0.682	ng/ul	99
19) Fluoranthene	19.518	202	56685	0.669	ng/ul#	94
21) Benzo(a)anthracene	21.622	228	59467	0.758	ng/ul	99
25) Benzo(k)fluoranthene	23.411	252	55936	0.741	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	26.728	276	49173	0.571	ng/ul#	99
29) Benzo(g,h,i)perylene	27.529	276	46237	0.633	ng/ul	96

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

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