

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092523\
 Data File : BM042051.D
 Acq On : 26 Sep 2023 02:25
 Operator : MA/JU
 Sample : 04429-01DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBKJ4DL

Quant Time: Sep 26 03:13:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:50:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	5361	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.763	136	12167	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.592	164	9215	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.346	188	12001	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.518	240	6462	0.400	ng/ul	0.00
23) Perylene-d12	23.924	264	6685	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	6463	0.885	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	1249	0.082	ng/ul	0.00
18) Fluoranthene-d10	19.366	212	2269	0.085	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.406	88	2642	0.344	ng/ul#	1
5) Naphthalene	10.812	128	38300	0.940	ng/ul	100
7) 2-Methylnaphthalene	12.418	142	20682	0.851	ng/ul	100
8) 1-Methylnaphthalene	12.638	142	12911	0.523	ng/ul	84
10) Acenaphthylene	14.342	152	1356	0.022	ng/ul#	1
11) Acenaphthene	14.624	153	1710	0.036	ng/ul#	50
12) Fluorene	15.642	166	2412	0.045	ng/ul#	80
14) Pentachlorophenol	16.978	266	46433	7.315	ng/ul	100

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

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