

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092523\
 Data File : BM042074.D
 Acq On : 26 Sep 2023 17:29
 Operator : MA/JU
 Sample : 04429-11DL 5X
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBBM3DL

Quant Time: Sep 27 04:26:27 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	6194	0.400	ng/ul	0.00
4) Naphthalene-d8	10.757	136	16202	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.592	164	8330	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	17185	0.400	ng/ul	-0.01
17) Chrysene-d12	21.516	240	8631	0.400	ng/ul	0.00
23) Perylene-d12	23.924	264	8272	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	6005	0.711	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.341	152	633	0.031	ng/ul	-0.01
18) Fluoranthene-d10	19.366	212	1082	0.030	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.406	88	3624	0.408	ng/ul	96
5) Naphthalene	10.807	128	161059	2.967	ng/ul	98
7) 2-Methylnaphthalene	12.412	142	50312	1.555	ng/ul	99
8) 1-Methylnaphthalene	12.632	142	32906	1.002	ng/ul	92
11) Acenaphthene	14.652	153	1653	0.039	ng/ul#	91
12) Fluorene	15.642	166	1236	0.026	ng/ul#	89
14) Pentachlorophenol	16.970	266	2980	0.328	ng/ul	100
15) Phenanthrene	17.384	178	2969	0.037	ng/ul#	85

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

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