

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100824\
 Data File : BF139705.D
 Acq On : 08 Oct 2024 11:43
 Operator : RC/JU
 Sample : P4290-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IB-4

Quant Time: Oct 08 12:06:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	89701	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	343039	20.000	ng	0.00
39) Acenaphthene-d10	9.898	164	194140	20.000	ng	-0.01
64) Phenanthrene-d10	11.386	188	373199	20.000	ng	0.00
76) Chrysene-d12	14.027	240	223398	20.000	ng	0.00
86) Perylene-d12	15.498	264	184671	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	712168	137.608	ng	0.02
7) Phenol-d6	6.504	99	883865	126.998	ng	0.00
23) Nitrobenzene-d5	7.428	82	662041	97.713	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	266778	142.349	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1173812	92.820	ng	0.00
79) Terphenyl-d14	12.975	244	1317872	96.796	ng	0.00
Target Compounds						
77) Benzidine	12.727	184	99517	25.255	ng	Qvalue 99

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

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