

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061824\
 Data File : BF138233.D
 Acq On : 18 Jun 2024 19:29
 Operator : CG\JU
 Sample : P2918-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 POLE

Quant Time: Jun 19 01:24:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 13 01:52:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	360979	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	1366590	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	700815	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	966545	20.000	ng	0.00
76) Chrysene-d12	14.039	240	627454	20.000	ng	-0.01
86) Perylene-d12	15.515	264	514810	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	2601104	120.539	ng	0.02
7) Phenol-d6	6.522	99	2958598	108.124	ng	0.00
23) Nitrobenzene-d5	7.457	82	2099655	89.445	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	764824	109.624	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	3790476	86.006	ng	0.00
79) Terphenyl-d14	12.998	244	3239493	81.772	ng	0.00
Target Compounds						
77) Benzidine	12.739	184	135841	18.717	ng	Qvalue 99

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

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