

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100423\
 Data File : BF135581.D
 Acq On : 04 Oct 2023 21:14
 Operator : CG\JU
 Sample : 04668-07MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200027MSD

Quant Time: Oct 04 23:28:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 04 17:06:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	150896	20.000	ng	0.00
21) Naphthalene-d8	8.016	136	589416	20.000	ng	0.00
39) Acenaphthene-d10	9.775	164	299073	20.000	ng	0.00
64) Phenanthrene-d10	11.263	188	588265	20.000	ng	# 0.00
76) Chrysene-d12	13.921	240	303995	20.000	ng	0.00
86) Perylene-d12	15.386	264	289105	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	1101118	118.685	ng	0.01
7) Phenol-d6	6.387	99	1299076	110.119	ng	0.00
23) Nitrobenzene-d5	7.304	82	918333	84.647	ng	0.00
42) 2,4,6-Tribromophenol	10.569	330	350464	117.919	ng	0.00
45) 2-Fluorobiphenyl	9.092	172	1575261	79.467	ng	0.00
79) Terphenyl-d14	12.863	244	1714200	87.414	ng	0.00

Target Compounds	Qvalue

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

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