

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040422\
 Data File : BF127632.D
 Acq On : 04 Apr 2022 15:15
 Operator : CG\JU
 Sample : N2249-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7D

Quant Time: Apr 04 16:26:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 29 07:13:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	77419	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	298450	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	173623	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	327765	20.000	ng	# 0.00
76) Chrysene-d12	14.016	240	289123	20.000	ng	0.00
86) Perylene-d12	15.504	264	208255	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	354680	73.128	ng	0.00
7) Phenol-d6	6.475	99	307681	46.409	ng	0.00
23) Nitrobenzene-d5	7.404	82	633655	90.756	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	233926	126.479	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	989097	83.187	ng	0.00
79) Terphenyl-d14	12.951	244	1259688	71.324	ng	0.00

Target Compounds	Qvalue

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

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