

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102522\
 Data File : BF130771.D
 Acq On : 25 Oct 2022 16:08
 Operator : CG\JU
 Sample : N5222-01DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12DL

Quant Time: Oct 26 13:48:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 21 06:12:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.563	152	145084	20.000	ng	0.00
21) Naphthalene-d8	7.851	136	447363	20.000	ng	0.00
39) Acenaphthene-d10	9.592	164	245879	20.000	ng	0.00
64) Phenanthrene-d10	11.075	188	421054	20.000	ng	0.00
76) Chrysene-d12	13.704	240	274554	20.000	ng	0.00
86) Perylene-d12	15.086	264	264700	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.193	112	106835	12.698	ng	0.03
7) Phenol-d6	6.246	99	83751	8.227	ng	0.02
23) Nitrobenzene-d5	7.134	82	148957	17.663	ng	0.00
42) 2,4,6-Tribromophenol	10.386	330	59273	25.349	ng	0.00
45) 2-Fluorobiphenyl	8.916	172	301766	18.703	ng	-0.01
79) Terphenyl-d14	12.651	244	298856	18.343	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	7.881	128	4301543	186.484	ng	96
37) 2-Methylnaphthalene	8.557	142	671519	44.432	ng	100
38) 1-Methylnaphthalene	8.657	142	1165944	77.812	ng	98
46) 1,1'-Biphenyl	9.016	154	215586	12.571	ng	97
52) Acenaphthene	9.622	154	201869	15.455	ng	100
58) Fluorene	10.139	166	90972	5.441	ng	100
71) Phenanthrene	11.098	178	178204	7.718	ng	100

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

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