

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102522\
 Data File : BF130772.D
 Acq On : 25 Oct 2022 16:42
 Operator : CG\JU
 Sample : N5222-01DL2 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12DL2

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 11/02/2022
 Supervised By :Jagrut Upadhyay 11/02/2022

Quant Time: Oct 26 15:01:30 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	116560	20.000	ng	0.00
21) Naphthalene-d8	7.845	136	450900	20.000	ng	0.00
39) Acenaphthene-d10	9.592	164	212978	20.000	ng	0.00
64) Phenanthrene-d10	11.074	188	329145	20.000	ng	0.00
76) Chrysene-d12	13.704	240	343124	20.000	ng	# 0.00
86) Perylene-d12	15.086	264	275101	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.210	112	47259	6.992	ng	0.05
7) Phenol-d6	6.263	99	32470m	3.970	ng	0.04
23) Nitrobenzene-d5	7.139	82	77020	9.061	ng	0.00
42) 2,4,6-Tribromophenol	10.386	330	21933	10.829	ng	0.00
45) 2-Fluorobiphenyl	8.916	172	136949	9.799	ng	-0.01
79) Terphenyl-d14	12.651	244	170330	8.365	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	7.875	128	2508308	107.889	ng	99
37) 2-Methylnaphthalene	8.557	142	335058	21.996	ng	100
38) 1-Methylnaphthalene	8.657	142	632424	41.875	ng	99
46) 1,1'-Biphenyl	9.022	154	91481	6.158	ng	98
52) Acenaphthene	9.622	154	79262	7.006	ng	98
58) Fluorene	10.139	166	33778	2.332	ng	99
71) Phenanthrene	11.098	178	70327	3.896	ng	99

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

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