

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042622\
 Data File : BF127936.D
 Acq On : 27 Apr 2022 02:27
 Operator : CG\JU
 Sample : N2567-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-14A-1A-4-2R

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 04/27/2022
 Supervised By :Jagrut Upadhyay 04/27/2022

Quant Time: Apr 27 03:21:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 00:43:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.939	152	126322	20.000	ng	0.00
21) Naphthalene-d8	8.228	136	455797	20.000	ng	# 0.00
39) Acenaphthene-d10	9.986	164	223669	20.000	ng	0.00
64) Phenanthrene-d10	11.480	188	316561	20.000	ng	0.00
76) Chrysene-d12	14.133	240	209217	20.000	ng	# 0.00
86) Perylene-d12	15.662	264	173487	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.569	112	805416	101.554	ng	0.00
7) Phenol-d6	6.569	99	1055416	102.558	ng	0.00
23) Nitrobenzene-d5	7.510	82	739814	76.326	ng	0.00
42) 2,4,6-Tribromophenol	10.780	330	212858	94.536	ng	0.00
45) 2-Fluorobiphenyl	9.304	172	1114971	84.332	ng	0.00
79) Terphenyl-d14	13.068	244	872313	76.240	ng	0.00
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.939	142	131304	8.558	ng	98
38) 1-Methylnaphthalene	9.039	142	148187	9.937	ng	100
71) Phenanthrene	11.504	178	67947m	4.100	ng	
78) Pyrene	12.927	202	46765	2.769	ng	# 82
83) Chrysene	14.157	228	35208	2.644	ng	# 56

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

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