

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041922\
 Data File : BF127852.D
 Acq On : 19 Apr 2022 20:37
 Operator : CG\JU
 Sample : N2452-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-15-1-4

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/20/2022
 Supervised By : Jagrut Upadhyay 04/20/2022

Quant Time: Apr 20 03:54:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF041922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 19 17:31:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.940	152	69942	20.000	ng	0.00
21) Naphthalene-d8	8.222	136	250524	20.000	ng	# 0.00
39) Acenaphthene-d10	9.986	164	149564	20.000	ng	0.00
64) Phenanthrene-d10	11.475	188	271318	20.000	ng	0.00
76) Chrysene-d12	14.121	240	177687	20.000	ng	#-0.01
86) Perylene-d12	15.639	264	142208	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	542772	121.575	ng	0.01
7) Phenol-d6	6.569	99	616774	110.628	ng	0.00
23) Nitrobenzene-d5	7.510	82	576285	106.682	ng	0.00
42) 2,4,6-Tribromophenol	10.780	330	262999	144.350	ng	0.00
45) 2-Fluorobiphenyl	9.304	172	1078027	94.145	ng	0.00
79) Terphenyl-d14	13.069	244	1155859	102.105	ng	0.00
Target Compounds						
32) Benzoic acid	7.928	122	6610m	2.663	ng	Qvalue

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

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