

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131139.D
 Acq On : 10 Nov 2022 23:44
 Operator : CG\JU
 Sample : N5494-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-20

Manual Integrations
 APPROVED

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

Quant Time: Nov 11 00:52:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	56627	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	216365	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	93056	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	156180	20.000	ng	0.00
76) Chrysene-d12	14.074	240	174673	20.000	ng	0.00
86) Perylene-d12	15.568	264	107283	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	363896	109.285	ng	0.00
7) Phenol-d6	6.516	99	424479	97.508	ng	0.00
23) Nitrobenzene-d5	7.451	82	274099	57.158	ng	-0.01
42) 2,4,6-Tribromophenol	10.721	330	79317	82.475	ng	-0.01
45) 2-Fluorobiphenyl	9.245	172	414459	62.175	ng	-0.01
79) Terphenyl-d14	13.010	244	420227	42.181	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.080	276	6188	4.513	ng	Qvalue # 83
88) Benzo(b)fluoranthene	15.127	252	15630m	2.061	ng	
92) Benzo(g,h,i)perylene	17.533	276	6821	5.188	ng	# 95

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

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