

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110422\
 Data File : BF131012.D
 Acq On : 05 Nov 2022 02:51
 Operator : CG\JU
 Sample : N5430-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 185465

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/05/2022
 Supervised By : Jagrut Upadhyay 11/07/2022

Quant Time: Nov 05 03:54:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 06:20:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.910	152	54427	20.000	ng	0.00
21) Naphthalene-d8	8.192	136	188447	20.000	ng	-0.01
39) Acenaphthene-d10	9.957	164	90884	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	157255	20.000	ng	-0.01
76) Chrysene-d12	14.098	240	180297	20.000	ng	0.00
86) Perylene-d12	15.604	264	191734	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	256934	81.839	ng	0.00
7) Phenol-d6	6.528	99	374632	91.826	ng	-0.01
23) Nitrobenzene-d5	7.475	82	283075	90.578	ng	-0.01
42) 2,4,6-Tribromophenol	10.745	330	52371	68.453	ng	-0.01
45) 2-Fluorobiphenyl	9.269	172	482713	80.301	ng	-0.01
79) Terphenyl-d14	13.039	244	540164	54.900	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.663	202	32385	3.755	ng	99
88) Benzo(b)fluoranthene	15.157	252	30510m	2.462	ng	
90) Benzo(a)pyrene	15.533	252	22442	2.252	ng	# 94

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

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