

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082522\
 Data File : BF130003.D
 Acq On : 25 Aug 2022 17:10
 Operator : CG\JU
 Sample : N4205-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-6

Quant Time: Aug 26 01:00:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 14:45:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.757	152	108972	20.000	ng	0.00
21) Naphthalene-d8	8.034	136	420429	20.000	ng	0.00
39) Acenaphthene-d10	9.786	164	204332	20.000	ng	0.00
64) Phenanthrene-d10	11.280	188	318589	20.000	ng	0.00
76) Chrysene-d12	13.927	240	230019	20.000	ng	0.00
86) Perylene-d12	15.374	264	246336	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	674019	102.409	ng	0.00
7) Phenol-d6	6.416	99	829339	100.623	ng	0.00
23) Nitrobenzene-d5	7.322	82	532165	67.759	ng	-0.01
42) 2,4,6-Tribromophenol	10.586	330	187876	91.612	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	944002	70.168	ng	0.00
79) Terphenyl-d14	12.863	244	807100	58.385	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.304	178	65293	3.670	ng	99
75) Fluoranthene	12.492	202	99452	5.462	ng	98
78) Pyrene	12.727	202	92182	4.612	ng	98
81) Benzo(a)anthracene	13.916	228	51348	3.245	ng	98
83) Chrysene	13.951	228	43232	2.919	ng	99
88) Benzo(b)fluoranthene	14.957	252	61804m	3.664	ng	
90) Benzo(a)pyrene	15.315	252	46219	3.479	ng	97

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

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