

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110822\
 Data File : BF131074.D
 Acq On : 08 Nov 2022 19:01
 Operator : CG\JU
 Sample : N5415-14RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-27RE

Quant Time: Nov 09 04:07:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 01:40:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	77166	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	280530	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	128928	20.000	ng	0.00
64) Phenanthrene-d10	11.439	188	187360	20.000	ng	0.00
76) Chrysene-d12	14.086	240	179919	20.000	ng	0.00
86) Perylene-d12	15.592	264	171089	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	508689	106.298	ng	0.00
7) Phenol-d6	6.528	99	621648	106.429	ng	0.00
23) Nitrobenzene-d5	7.463	82	406394	67.150	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	118013	83.139	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	631426	66.744	ng	0.00
79) Terphenyl-d14	13.027	244	338512	30.165	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.098	276	4804	3.633	ng	Qvalue # 87
92) Benzo(g,h,i)perylene	17.562	276	4085	3.708	ng	# 91

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

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