

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110822\
 Data File : BF131073.D
 Acq On : 08 Nov 2022 18:29
 Operator : CG\JU
 Sample : N5415-13RE
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-26RE

Quant Time: Nov 09 04:07:45 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	66554	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	229579	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	108158	20.000	ng	0.00
64) Phenanthrene-d10	11.439	188	202607	20.000	ng	0.00
76) Chrysene-d12	14.086	240	183727	20.000	ng	0.00
86) Perylene-d12	15.586	264	181492	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	389713	94.421	ng	0.00
7) Phenol-d6	6.522	99	443665	88.069	ng	0.00
23) Nitrobenzene-d5	7.463	82	280283	56.590	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	89984	75.567	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	411576	51.860	ng	0.00
79) Terphenyl-d14	13.027	244	482318	42.088	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.103	276	5215	3.640	ng	Qvalue # 83
92) Benzo(g,h,i)perylene	17.568	276	5152	3.768	ng	# 89

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

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