

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031723\
 Data File : BP014134.D
 Acq On : 17 Mar 2023 22:47
 Operator : CG/JU
 Sample : 01936-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OR-03-03152023

Quant Time: Mar 18 00:29:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 18 00:19:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	140235	20.000	ng	0.00
21) Naphthalene-d8	10.887	136	539735	20.000	ng	0.00
39) Acenaphthene-d10	14.704	164	343966	20.000	ng	0.00
64) Phenanthrene-d10	17.463	188	761609	20.000	ng	0.00
76) Chrysene-d12	21.545	240	716921	20.000	ng	0.00
86) Perylene-d12	24.092	264	851449	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.605	112	897332	103.772	ng	0.00
7) Phenol-d6	7.222	99	1173482	103.362	ng	0.00
23) Nitrobenzene-d5	9.234	82	777983	65.881	ng	0.00
42) 2,4,6-Tribromophenol	16.198	330	575796	103.247	ng	0.00
45) 2-Fluorobiphenyl	13.328	172	1635330	61.596	ng	0.00
79) Terphenyl-d14	20.004	244	2538991	58.398	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.504	178	95178	2.237	ng	98
75) Fluoranthene	19.463	202	182766	3.594	ng	97
78) Pyrene	19.816	202	182984	3.325	ng	98
83) Chrysene	21.586	228	104791	2.104	ng #	82
88) Benzo(b)fluoranthene	23.310	252	119961	2.152	ng #	46

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

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