

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080422\
 Data File : BM036102.D
 Acq On : 04 Aug 2022 17:14
 Operator : CG/JU
 Sample : N3931-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-117

Quant Time: Aug 05 03:39:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	355387	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	1593964	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	998611	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	2073253	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1567240	20.000	ng	0.00
86) Perylene-d12	23.791	264	1316674	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	2001946	91.328	ng	0.00
7) Phenol-d6	7.034	99	2650150	95.014	ng	0.00
23) Nitrobenzene-d5	9.028	82	1589988	55.837	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1094383	93.417	ng	0.00
45) 2-Fluorobiphenyl	13.127	172	3374613	49.895	ng	0.00
79) Terphenyl-d14	19.868	244	4368975	54.986	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.280	178	549030	5.198	ng	100
75) Fluoranthene	19.298	202	953849	8.067	ng	98
78) Pyrene	19.662	202	841291	8.657	ng	100
81) Benzo(a)anthracene	21.409	228	410801	4.256	ng	98
83) Chrysene	21.462	228	404023	4.403	ng	98
88) Benzo(b)fluoranthene	23.068	252	470235	6.096	ng	# 94
90) Benzo(a)pyrene	23.680	252	279211	4.314	ng	# 93

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

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