

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011022\
 Data File : BM033689.D
 Acq On : 10 Jan 2022 16:40
 Operator : CG/JU
 Sample : N1081-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OB-CS-12

Quant Time: Jan 11 02:25:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 05 17:07:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.978	152	211268	20.000	ng	0.00
21) Naphthalene-d8	10.795	136	869403	20.000	ng	#-0.01
39) Acenaphthene-d10	14.613	164	518687	20.000	ng	-0.01
64) Phenanthrene-d10	17.348	188	913308	20.000	ng	-0.01
76) Chrysene-d12	21.518	240	660535	20.000	ng	0.00
86) Perylene-d12	23.947	264	711843	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	212596	15.871	ng	0.00
7) Phenol-d6	7.131	99	466764	25.854	ng	0.00
23) Nitrobenzene-d5	9.142	82	1173028	62.311	ng	-0.01
42) 2,4,6-Tribromophenol	16.095	330	76628	14.144	ng	-0.01
45) 2-Fluorobiphenyl	13.248	172	2553290	64.476	ng	-0.01
79) Terphenyl-d14	19.965	244	2624047	71.507	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.389	178	117469	2.248	ng	100
78) Pyrene	19.759	202	964567	20.741	ng	98
81) Benzo(a)anthracene	21.500	228	193826	4.327	ng	# 61
83) Chrysene	21.553	228	244840	5.712	ng	# 88
90) Benzo(a)pyrene	23.835	252	183981	4.884	ng	# 84
92) Benzo(g,h,i)perylene	27.265	276	239058	5.577	ng	# 92

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

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