

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060321\
 Data File : BP005699.D
 Acq On : 03 Jun 2021 22:04
 Operator : CG/JU
 Sample : M2580-02 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B1(2-4)

Quant Time: Jun 04 01:05:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 27 16:11:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	195766	20.00	ng	0.00
21) Naphthalene-d8	10.793	136	758610	20.00	ng	# 0.00
39) Acenaphthene-d10	14.610	164	465987	20.00	ng	0.00
64) Phenanthrene-d10	17.351	188	964552	20.00	ng	# 0.00
76) Chrysene-d12	21.422	240	956234	20.00	ng	# 0.00
86) Perylene-d12	23.863	264	1054288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	89093	7.08	ng	0.00
7) Phenol-d6	7.146	99	114677	6.68	ng	0.00
23) Nitrobenzene-d5	9.146	82	69903	5.12	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	41728	6.98	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	171471	4.87	ng	0.00
79) Terphenyl-d14	19.892	244	260077	5.10	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.392	178	181233	3.02	ng	98
75) Fluoranthene	19.351	202	316596	4.39	ng	95
78) Pyrene	19.704	202	283281	4.20	ng	99
81) Benzo(a)anthracene	21.404	228	170275	2.43	ng	# 94
83) Chrysene	21.457	228	146658	2.16	ng	# 95
88) Benzo(b)fluoranthene	23.116	252	199281	2.61	ng	# 79
90) Benzo(a)pyrene	23.751	252	159604	2.29	ng	# 81

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

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