

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100821\
 Data File : BP007385.D
 Acq On : 08 Oct 2021 21:34
 Operator : CG/JU
 Sample : M4112-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CON-WALL

Quant Time: Oct 09 06:19:23 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP100821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 08 15:35:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	202124	20.00	ng	0.00
21) Naphthalene-d8	10.575	136	443769	20.00	ng	0.00
39) Acenaphthene-d10	14.422	164	300815	20.00	ng	0.00
64) Phenanthrene-d10	17.175	188	651625	20.00	ng	-0.01
76) Chrysene-d12	21.269	240	581666	20.00	ng	-0.01
86) Perylene-d12	23.621	264	573114	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	745278	60.76	ng	0.00
7) Phenol-d6	6.963	99	1149193	69.78	ng	0.00
23) Nitrobenzene-d5	8.946	82	338747	44.96	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	392887	79.62	ng	-0.01
45) 2-Fluorobiphenyl	13.045	172	988071	45.90	ng	0.00
79) Terphenyl-d14	19.745	244	1612673	48.90	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.216	178	375981	10.93	ng	99
75) Fluoranthene	19.192	202	514240	11.59	ng	98
78) Pyrene	19.545	202	362241	9.04	ng	99
81) Benzo(a)anthracene	21.251	228	120857	3.22	ng	99
83) Chrysene	21.304	228	172808	4.89	ng	98
87) Indeno(1,2,3-cd)pyrene	26.074	276	88310	2.25	ng	# 94
88) Benzo(b)fluoranthene	22.904	252	192711	2.69	ng	96
90) Benzo(a)pyrene	23.516	252	112883	3.90	ng	99
92) Benzo(g,h,i)perylene	26.827	276	111540	3.63	ng	# 96

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

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