

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005267.D
 Acq On : 12 Apr 2021 15:43
 Operator : CG/JU
 Sample : M1992-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AU-05-040921

Quant Time: Apr 12 16:56:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	29455	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	100585	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	59806	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	133457	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	166659	20.00	ng	0.00
86) Perylene-d12	23.480	264	172294	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	289841	103.27	ng	0.00
7) Phenol-d6	6.863	99	319612	99.80	ng	0.00
23) Nitrobenzene-d5	8.840	82	210312	65.94	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	72688	96.28	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	289183	65.28	ng	0.00
79) Terphenyl-d14	19.680	244	498939	58.81	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.127	202	39712	4.29	ng	98
78) Pyrene	19.480	202	38779	3.74	ng	98
81) Benzo(a)anthracene	21.186	228	33912	3.08	ng	98
83) Chrysene	21.239	228	34535	3.23	ng	97
87) Indeno(1,2,3-cd)pyrene	25.798	276	29894	2.27	ng	# 92
88) Benzo(b)fluoranthene	22.792	252	54871	4.87	ng	98
90) Benzo(a)pyrene	23.374	252	39016	3.74	ng	97
92) Benzo(g,h,i)perylene	26.515	276	34042	3.07	ng	94

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

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