

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005274.D
 Acq On : 12 Apr 2021 19:41
 Operator : CG/JU
 Sample : M1997-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-58-56

Quant Time: Apr 13 00:27:11 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	7.675	152	36087	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	124194	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	77257	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	171834	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	181612	20.00	ng	0.00
86) Perylene-d12	23.474	264	183079	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	351732	102.29	ng	0.00
7) Phenol-d6	6.863	99	389076	99.17	ng	0.00
23) Nitrobenzene-d5	8.840	82	259092	65.79	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	102444	105.05	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	362053	63.27	ng	0.00
79) Terphenyl-d14	19.680	244	588281	63.63	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.145	178	19661	2.02	ng	# 79
75) Fluoranthene	19.127	202	67235	5.64	ng	95
78) Pyrene	19.480	202	63392	5.62	ng	96
81) Benzo(a)anthracene	21.186	228	33782	2.82	ng	98
83) Chrysene	21.239	228	40269	3.45	ng	99
87) Indeno(1,2,3-cd)pyrene	25.798	276	29343	2.10	ng	# 91
88) Benzo(b)fluoranthene	22.786	252	58598	4.89	ng	99
90) Benzo(a)pyrene	23.374	252	37425	3.37	ng	95
92) Benzo(g,h,i)perylene	26.515	276	32851	2.78	ng	# 90

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

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